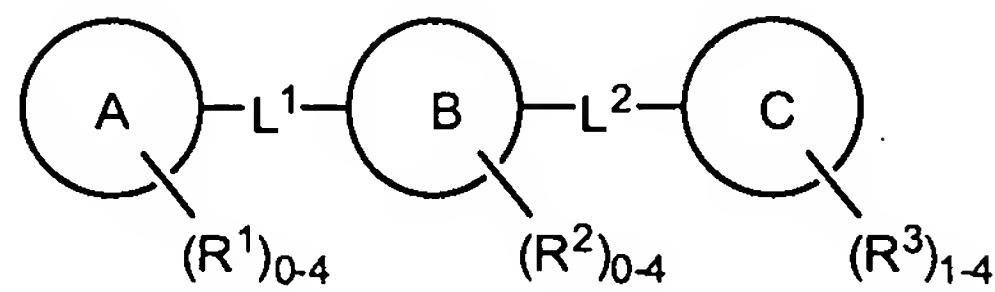


The Listing of the Claims:

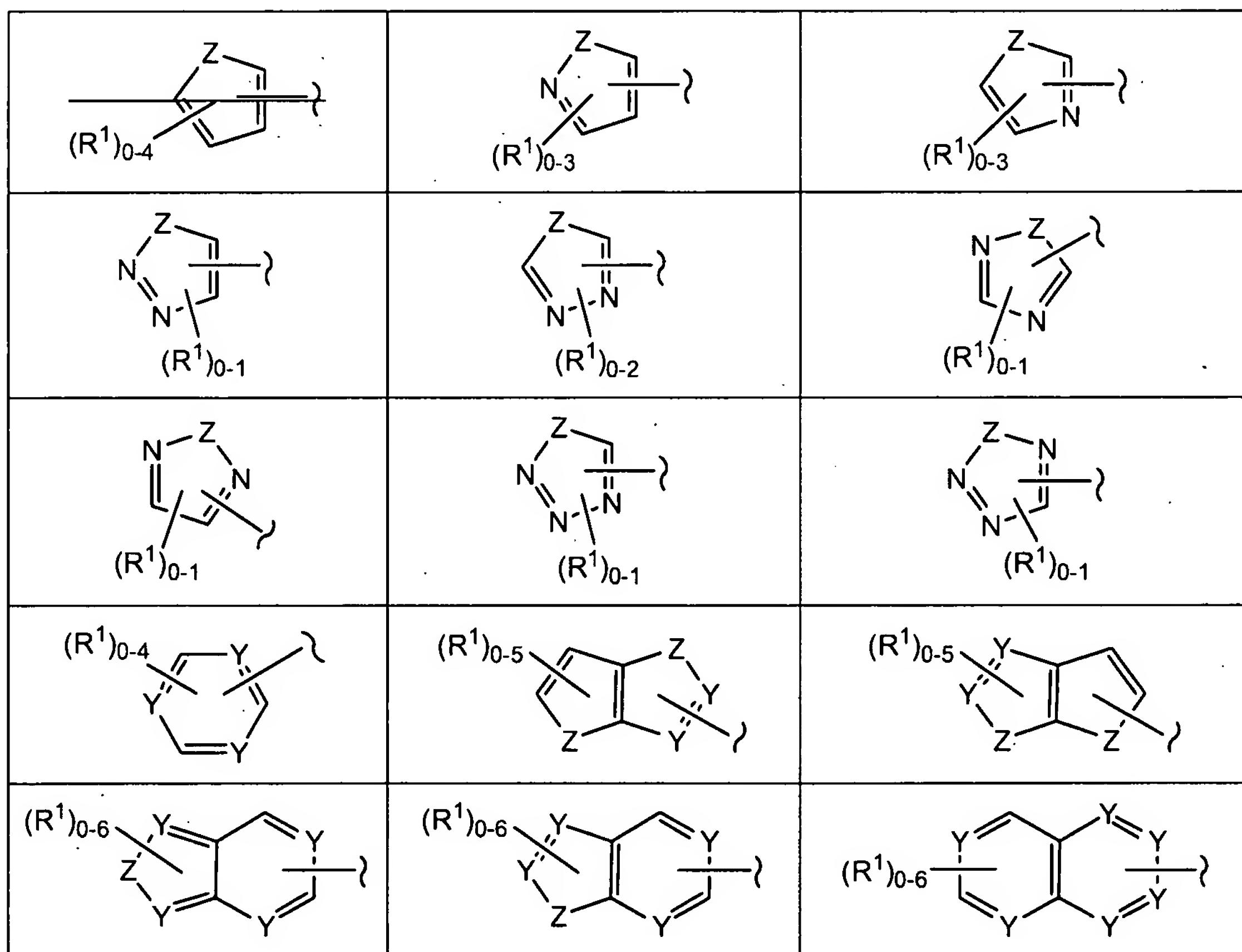
1. (currently amended) A compound for modulating c-Kit activity according to Formula I,



I

or a pharmaceutically acceptable salt, thereof, wherein,

ring A is:



wherein each Y is independently either =C(H)- or =N-; and Z is selected from -O-, -S-, and -N(R⁷)-;

each R¹ is independently selected from -H, halogen, trihalomethyl, -CN, -NO₂, -OR⁴, -S(O)₀₋₂R⁴, -SO₂N(R⁴)R⁴, -C(=O)N(R⁴)R⁴, -C(=NR⁵)N(R⁴)R⁴, -C(=NR⁵)R⁴, -N(R⁴)SO₂R⁴, -N(R⁴)C(O)R⁴, -NCO₂R⁴, -C(=O)R⁴, optionally substituted alkoxy, optionally substituted C₁₋₆alkyl, optionally substituted aryl C₁₋₆alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C₁₋₆alkyl;

two adjacent of R¹, together with the annular atoms to which they are attached, can form a five- to six-membered ring containing up to two heteroatoms and optionally substituted with up to three of R¹⁰;

L¹ is a single bond;

ring B is phenyl a five- to ten-membered aryl or a five- to ten-membered heterocyclyl;

each R² is independently selected from -H, halogen, trihalomethyl, -CN, -NO₂, -OR⁴, -N(R⁴)R⁴, -S(O)₀₋₂R⁴, -SO₂N(R⁴)R⁴, -C(=O)N(R⁴)R⁴, -C(=NR⁵)N(R⁴)R⁴, -C(=NR⁵)R⁴, -N(R⁴)SO₂R⁴, -N(R⁴)C(O)R⁴, -NCO₂R⁴, -C(=O)R⁴, optionally substituted alkoxy, optionally substituted C₁₋₆alkyl, optionally substituted aryl, optionally substituted aryl C₁₋₆alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C₁₋₆alkyl;

two adjacent of R², together with the annular atoms to which they are attached, can form a five- to six-membered ring containing up to two heteroatoms and optionally substituted with up to three of R¹⁵;

L² is selected from -N(H)N(H)C(=O)N(H)-, -CH₂N(H)C(=O)N(H)-, -CH₂OC(=O)N(H)-, and -XCH₂C(=O)N(H)-; wherein X is selected from -O-, -S(O)₀₋₂-, and -N(R⁷)-; and any C-H of L² is optionally C-R²⁰;

ring C is phenyl or pyridyl;

each R³ is independently selected from halogen, trihalomethyl, -CN,

-NO_2 , -OR^4 , $\text{-N(R}^4\text{)R}^4$, $\text{-S(O)}_{0-2}\text{R}^4$, $\text{-SO}_2\text{N(R}^4\text{)R}^4$, $\text{-C(=O)N(R}^4\text{)R}^4$, $\text{-C(=NR}^5\text{)N(R}^4\text{)R}^4$, $\text{-C(=NR}^5\text{)R}^4$, $\text{-N(R}^4\text{)SO}_2\text{R}^4$, $\text{-N(R}^4\text{)C(O)R}^4$, $\text{-NCO}_2\text{R}^4$, -C(=O)R^4 , optionally substituted alkoxy, optionally substituted $\text{C}_{1-6}\text{alkyl}$, optionally substituted aryl, optionally substituted aryl $\text{C}_{1-6}\text{alkyl}$, optionally substituted heterocyclyl, and optionally substituted heterocyclyl $\text{C}_{1-6}\text{alkyl}$; provided R^3 is not a cyclic sulfonamide attached to ring C via the nitrogen of said cyclic sulfonamide,

wherein there exists at least one of R^3 that is halogen or trihalomethyl;

~~two adjacent of R^3 , together with the annular atoms to which they are attached, can form a five- to six-membered ring containing up to two heteroatoms and optionally substituted with up to three of R^{25} ;~~

R^4 is selected from -H, optionally substituted $\text{C}_{1-6}\text{alkyl}$, optionally substituted aryl, optionally substituted aryl $\text{C}_{1-6}\text{alkyl}$, optionally substituted heterocyclyl, and optionally substituted heterocyclyl $\text{C}_{1-6}\text{alkyl}$;

two of R^4 , when taken together with a common nitrogen to which they are attached, form an optionally substituted five- to seven-membered heterocyclyl, said optionally substituted five- to seven-membered heterocyclyl optionally containing at least one additional heteroatom selected from N, O, S, and P;

R^5 is selected from -H, -CN, -NO_2 , -OR^4 , $\text{-S(O)}_{0-2}\text{R}^4$, $\text{-CO}_2\text{R}^4$, optionally substituted $\text{C}_{1-6}\text{alkyl}$, optionally substituted $\text{C}_{1-6}\text{alkenyl}$, and optionally substituted $\text{C}_{1-6}\text{alkynyl}$;

R^7 is selected from -H, optionally substituted $\text{C}_{1-6}\text{alkyl}$, $\text{-SO}_2\text{N(R}^4\text{)R}^4$, $\text{-CO}_2\text{R}^4$, $\text{-C(=O)N(R}^4\text{)R}^4$, $\text{-C(=NR}^5\text{)N(R}^4\text{)R}^4$, $\text{-C(=NR}^5\text{)R}^4$, -C(=O)R^4 , optionally substituted alkoxy, optionally substituted aryl $\text{C}_{1-6}\text{alkyl}$, optionally substituted heterocyclyl, and optionally substituted heterocyclyl $\text{C}_{1-6}\text{alkyl}$; and

each of R^{10} , each of R^{15} , each of R^{20} , and each of R^{25} is independently selected from -H, halogen, trihalomethyl, -CN, -NO_2 , -OR^4 , $\text{-N(R}^4\text{)R}^4$, $\text{-S(O)}_{0-2}\text{R}^4$, $\text{-SO}_2\text{N(R}^4\text{)R}^4$, $\text{-CO}_2\text{R}^4$, $\text{-C(=O)N(R}^4\text{)R}^4$, $\text{-C(=NR}^5\text{)N(R}^4\text{)R}^4$, $\text{-C(=NR}^5\text{)R}^4$, $\text{-N(R}^4\text{)SO}_2\text{R}^4$, $\text{-N(R}^4\text{)C(O)R}^4$, $\text{-NCO}_2\text{R}^4$, -C(=O)R^4 , optionally substituted alkoxy, optionally substituted $\text{C}_{1-6}\text{alkyl}$,

optionally substituted aryl, optionally substituted aryl C₁₋₆alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C₁₋₆alkyl;

provided:

1) when both ring B and ring C are phenyl:

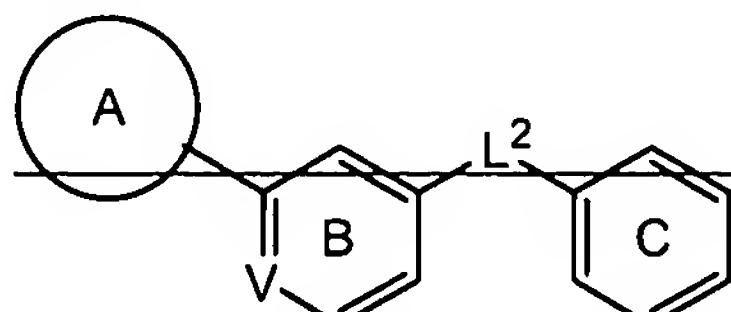
- a) and the compound comprises ring B CH₂N(H)C(=O)N(H) ring C, then L⁺ must be a single bond; R³ can not comprise a group of the formula -O(CH₂)₂₋₄N-piperazine that is *ortho* to L²; and ring A cannot be a 5-methyl [1,2,4] oxadiazol 3-yl radical, a 4H [1,2,4] oxadiazol 5-one 3-yl radical, nor a 4' [2,2';6',2'']terpyridinyl radical;
- b) and L⁺ is single bond, then L² cannot comprise N(H)C(=O)C(=O)N(H) nor -N(H)C(=Q)C(H)CNC(=O) (where Q is S or O);
- c) and L⁺ is other than single bond, then A cannot be quinolin 2-yl L⁺, quinolin 3-yl L⁺, or quinolin 4-yl L⁺;

2) when ring A is a fused aryl system, then L⁺ must be a single bond;

3) when ring B is phenyl, ring C is a C₆₋₁₆carboyclic, L⁺ is a single bond, and the compound comprises ring B OCH₂C(=O)N(H) then ring A cannot be a 2,5-dimethyl 1H-pyrrole 1-yl radical;

4) ring A cannot be a pyrimidin 2-yl radical when L⁺ is N(H) and ring B is phenyl;

5) when the compound comprises the formula,



where V is -C(H) or -N, and there is a nitrogen of L² bound directly to ring B, then A can not comprise a [1,2,4] oxadiazol 3-yl radical; and

6) the compound is not one of: N-naphthalen 1-yl 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide, N-[4-(phenyloxy)phenyl] 2-[[3-(1H-tetrazol-1-

yl)phenyl]oxy}acetamide, ~~N-(3,4-dimethylphenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide,~~ ~~N-(2,3-dimethylphenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide,~~ ~~N-(2,4-dimethylphenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide,~~ ~~N-(2,5-dimethylphenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide,~~ ~~N-(3,5-dimethylphenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide,~~ ~~N-(2,6-dimethylphenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide,~~ ~~2-{{3-(1H-tetrazol-1-yl)phenyl]oxy}-N-(2,4,6-trimethylphenyl)acetamide,~~ ~~N-(2-ethylphenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide,~~ ~~N-(4-ethylphenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide,~~ ~~N-(2,6-diethylphenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide,~~ ~~N-[2-(methyloxy)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide,~~ ~~N-[2-(ethyloxy)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide,~~ ~~N-[3-(ethyloxy)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide,~~ ~~N-[2,4-bis(methyloxy)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide,~~ ~~N-[4-(dimethylamino)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide,~~ ~~N-(2,3-dichlorophenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide,~~ ~~N-(4-chloro-3-methylphenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide,~~ ~~N-(4-bromophenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide,~~ ~~N-(2-fluorophenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide,~~ ~~N-(4-fluorophenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide,~~ ~~2-{{3-(1H-tetrazol-1-yl)phenyl]oxy}-N-[2-(trifluoromethyl)phenyl]acetamide,~~ ~~2-{{3-(1H-tetrazol-1-yl)phenyl]oxy}-N-[3-(trifluoromethyl)phenyl]acetamide,~~ ~~methyl-4-{{3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl]amino]benzoate,~~ ~~ethyl-4-{{3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl]amino]benzoate,~~ ~~3-{{3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl]amino]benzoic acid,~~ ~~N-[3-(methyloxy)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide,~~ ~~N-[4-(methyloxy)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide,~~ ~~N-[2-chloro-5-(trifluoromethyl)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide,~~ ~~N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide,~~

{[3-(4H-1,2,4-triazol-4-yl)phenyl]oxy} acetamide or acetamide, N-(4-chlorophenyl)-2- {[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(4-aminophenyl)-2 {[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, and N-(4-acetylphenyl)-2 {[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide.

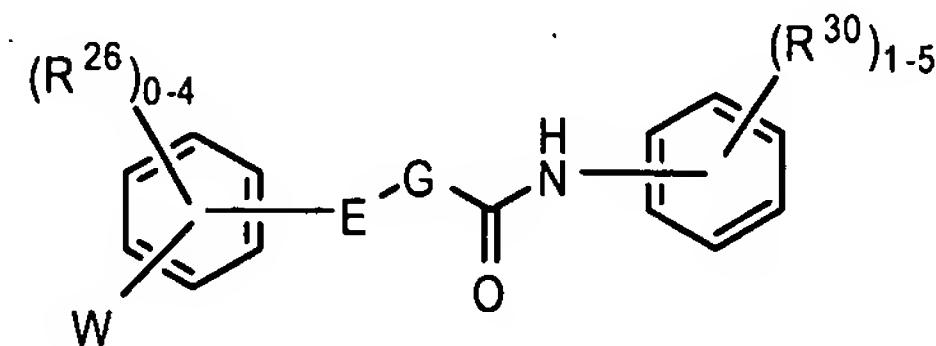
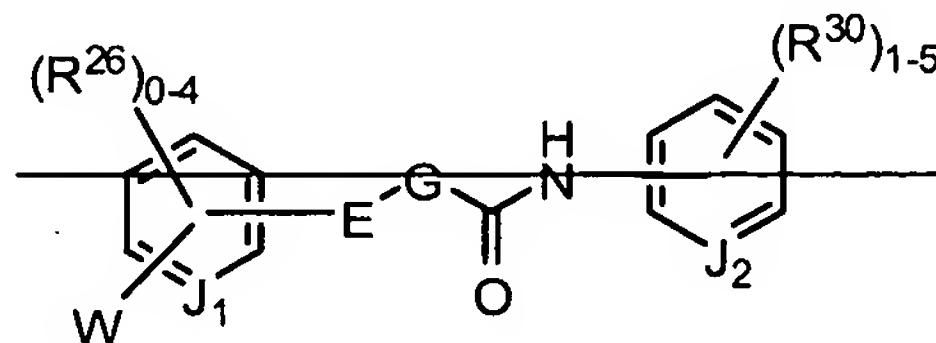
2-12. (cancelled)

13. (Currently Amended) The compound according to claim 10, 1, wherein there exists at least one of R³ that is trifluoromethyl.

14. (original) The compound according to claim 13, wherein ring C is a phenyl comprising a trifluoromethyl radical *meta*- to L².

15. (currently amended) The compound according to claim 1, claim 10, wherein each of R³ is independently selected from halogen, trihalomethyl, -OR⁴, -C(=O)R⁴, and optionally substituted C₁₋₆alkyl.

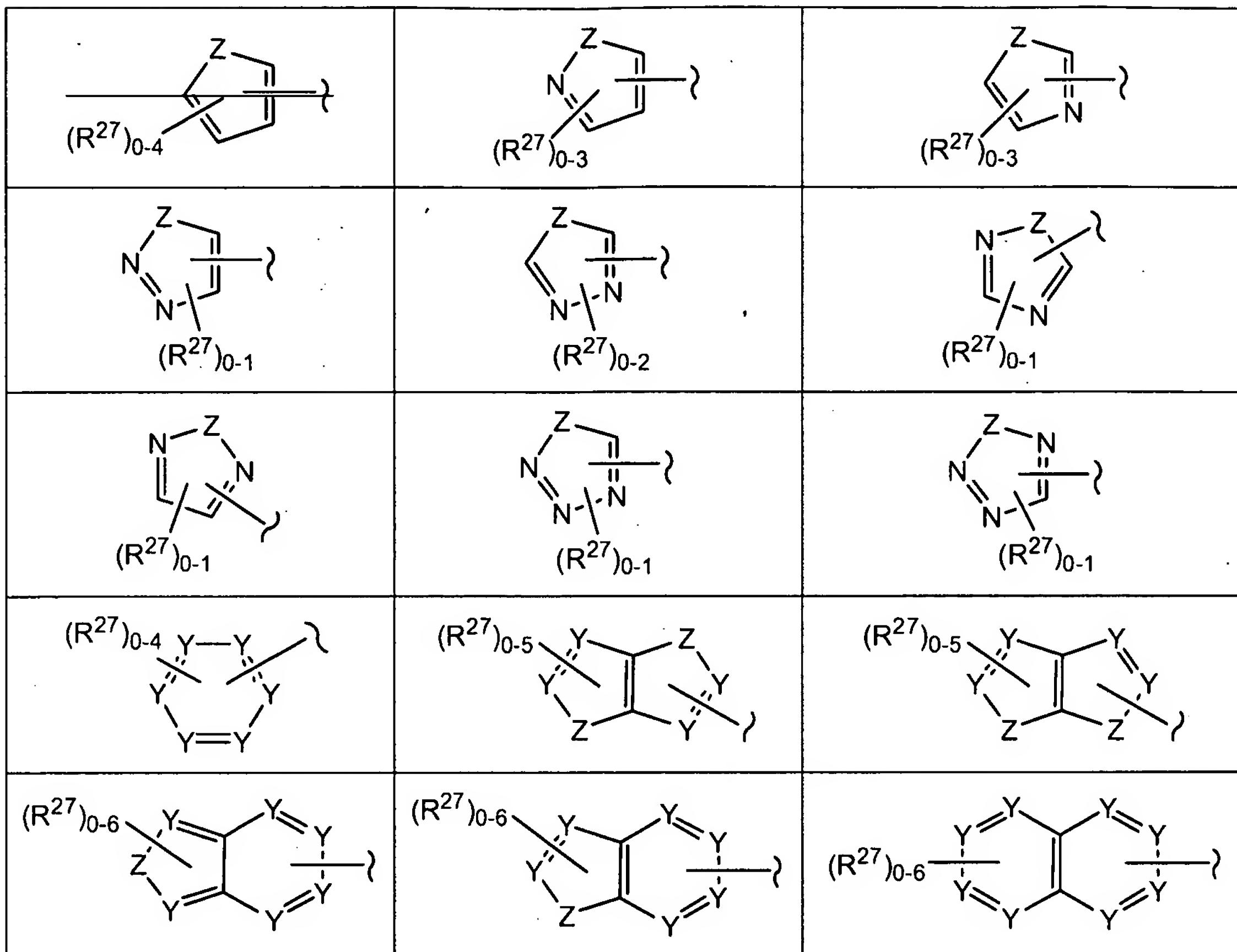
16. (currently amended) A compound for modulating c-Kit activity according to the following Formula: II,



II

or a pharmaceutically acceptable salt, thereof, wherein,

W is selected from the following:



each of R²⁷ independently selected from halogen, trihalomethyl, -CN, -NO₂, -OR⁵⁵, -S(O)₀₋₂R⁵⁵, -SO₂N(R⁵⁵)R⁵⁵, -C(=O)N(R⁵⁵)R⁵⁵, -C(=NR⁵⁰)N(R⁵⁵)R⁵⁵, -C(=NR⁵⁰)R⁵⁵, -N(R⁵⁵)SO₂R⁵⁵, -N(R⁵⁵)C(O)R⁵⁵, -NCO₂R⁵⁵, -C(=O)R⁵⁵, optionally substituted alkoxy, optionally substituted C₁₋₆alkyl, optionally substituted aryl C₁₋₆alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C₁₋₆alkyl;

each Y is independently either =C(H)- or =N-;

Z is selected from -O-, -S(O)₀₋₂-, and -N(R⁷)-

E and G are each independently selected from -O-, -S(O)₀₋₂-, -C(R³¹)R³²-, and -N(R³³)-;

J₁ and J₂ are each independently =C(H)- or =N-;

R^{26} is independently selected from -H, halogen, trihalomethyl, -CN, -NO₂, -OR⁴⁰, -N(R⁴⁰)R⁴⁰, -S(O)₀₋₂R⁴⁰, -SO₂N(R⁴⁰)R⁴⁰, -CO₂R⁴⁰, -C(=O)N(R⁴⁰)R⁴⁰, -C(=NR⁵⁰)N(R⁴⁰)R⁴⁰, -C(=NR⁵⁰)R⁴⁰, -N(R⁴⁰)SO₂R⁴⁰, -N(R⁴⁰)C(O)R⁴⁰, -NCO₂R⁴⁰, -C(=O)R⁴⁰, optionally substituted alkoxy, optionally substituted C₁₋₆alkyl, optionally substituted aryl, optionally substituted aryl C₁₋₆alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C₁₋₆alkyl;

R^{30} is independently selected from halogen, trihalomethyl, -CN, -NO₂, -OR⁴⁰, -N(R⁴⁰)R⁴⁰, -S(O)₀₋₂R⁴⁰, -SO₂N(R⁴⁰)R⁴⁰, -C(=O)N(R⁴⁰)R⁴⁰, -C(=NR⁵⁰)N(R⁴⁰)R⁴⁰, -C(=NR⁵⁰)R⁴⁰, -N(R⁴⁰)SO₂R⁴⁰, -N(R⁴⁰)C(O)R⁴⁰, -NCO₂R⁴⁰, -C(=O)R⁴⁰, optionally substituted alkoxy, optionally substituted C₁₋₆alkyl, optionally substituted aryl, optionally substituted aryl C₁₋₆alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C₁₋₆alkyl, wherein there exists at least one of R^{30} that is trihalomethyl; or

~~two adjacent of R^{26} or two adjacent of R^{30} , together with the annular atoms to which they are attached, can form a five to six membered ring containing up to two heteroatoms and optionally substituted with up to three of R^{35} ;~~

R^{31} and R^{32} are each independently selected from -H, halogen, trihalomethyl, -CN, -NO₂, -OR⁴⁰, -N(R⁴⁰)R⁴⁰, -S(O)₀₋₂R⁴⁰, -SO₂N(R⁴⁰)R⁴⁰, -CO₂R⁴⁰, -C(=O)N(R⁴⁰)R⁴⁰, -C(=NR⁵⁰)N(R⁴⁰)R⁴⁰, -C(=NR⁵⁰)R⁴⁰, -N(R⁴⁰)SO₂R⁴⁰, -N(R⁴⁰)C(O)R⁴⁰, -NCO₂R⁴⁰, -C(=O)R⁴⁰, optionally substituted alkoxy, optionally substituted C₁₋₆alkyl, optionally substituted aryl, optionally substituted aryl C₁₋₆alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C₁₋₆alkyl;

R^{33} is selected from -H, optionally substituted lower alkyl, -SO₂N(R⁴⁰)R⁴⁰, -CO₂R⁴⁰, -C(=O)N(R⁴⁰)R⁴⁰, -C(=NR⁵⁰)N(R⁴⁰)R⁴⁰, -C(=NR⁵⁰)R⁴⁰, -C(=O)R⁴⁰, optionally substituted alkoxy, optionally substituted C₁₋₆alkyl, optionally substituted aryl, optionally substituted aryl C₁₋₆alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C₁₋₆alkyl;

R^{40} is selected from -H, optionally substituted alkoxy, optionally substituted C₁₋₆alkyl, optionally substituted aryl, optionally substituted aryl C₁₋₆alkyl, optionally substituted heterocyclyl,

and optionally substituted heterocyclyl C₁₋₆alkyl;

two of R⁴⁰, when taken together with a common nitrogen to which they are attached, form an optionally substituted five- to seven-membered heterocyclyl, said optionally substituted five- to seven-membered heterocyclyl optionally containing at least one additional heteroatom selected from N, O, S, and P;

R⁵⁰ is selected from -H, -CN, -NO₂, -OR⁴⁰, -S(O)₀₋₂R⁴⁰, -CO₂R⁴⁰, optionally substituted C₁₋₆alkyl, optionally substituted C₁₋₆alkenyl, and optionally substituted C₁₋₆alkynyl;

R⁵⁵ is selected from -H, optionally substituted C₁₋₆alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C₁₋₆alkyl; and

two of R⁵⁵, when taken together with a common nitrogen to which they are attached, form an optionally substituted five- to seven-membered heterocyclyl, said optionally substituted five- to seven-membered heterocyclyl optionally containing at least one additional heteroatom selected from N, O, S, and P.

17. (cancelled)

18. (currently amended) The compound according to claim 17 16, wherein R³⁰ is selected from halogen, trihalomethyl, -OR⁴⁰, -N(R⁴⁰)R⁴⁰, -C(=O)R⁴⁰, optionally substituted alkoxy, optionally substituted C₁₋₆alkyl, optionally substituted aryl, optionally substituted aryl C₁₋₆alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C₁₋₆alkyl, wherein there exists at least one of R³⁰ that is trifluoromethyl.

19. (cancelled)

20. (cancelled)

21. (cancelled)

22. (cancelled)

23. (withdrawn from consideration, currently amended) The compound according to claim 22, claim 16, wherein E is selected from -O-, -S(O)₀₋₂-, and -NH-; and G is -CH₂-.

24. (withdrawn from consideration, currently amended) The compound according to claim 22, claim 16, wherein E is either -CH₂- or -NH-; and G is selected from -O-, -S-, and -NH-.

25. (cancelled)

26. (cancelled)

27. (currently amended) ~~The compound according to claim 1, selected from Table 3:~~

A compound selected from the following Table:

Table 3

Entry	Name	Structure
1	N-[5-chloro-2-(methyloxy)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}acetamide	
2	N-phenyl-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}acetamide	
3	N-(2-methylphenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}acetamide	

Table 3

Entry	Name	Structure
4	N-(2-chlorophenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}acetamide	
5	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}acetamide	
6	ethyl 2-{{{{3-(1H-tetrazol-1-yl)phenyl}oxy}acetyl}amino}4,5,6,7-tetrahydro-1-benzothiophene-3-carboxylate	
7	N-(3-chloro-2-methylphenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}acetamide	
8	N-(3-fluorophenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}acetamide	

Table 3

Entry	Name	Structure
9	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[3-(2H-tetrazol-5- yl)phenyl]oxy}acetamide	
10	N-(4-chloro-2-fluorophenyl)-2-{{3-(1H- tetrazol-1-yl)phenyl}oxy}acetamide	
11	N-(4-bromo-3-methylphenyl)-2-{{3-(1H- tetrazol-1-yl)phenyl}oxy}acetamide	
12	N-(4-morpholin-4-ylphenyl)-2-{{3-(1H- tetrazol-1-yl)phenyl}oxy}acetamide	
13	N-[4-fluoro-3-(trifluoromethyl)phenyl]-2- {[3-(1H-tetrazol-1- yl)phenyl]oxy}acetamide	

Table 3

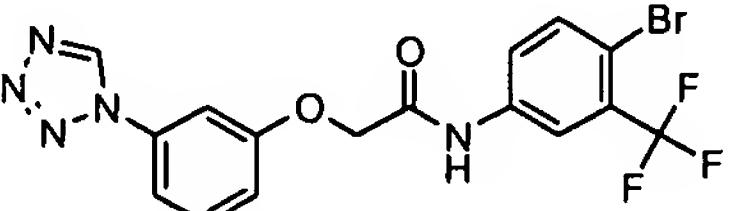
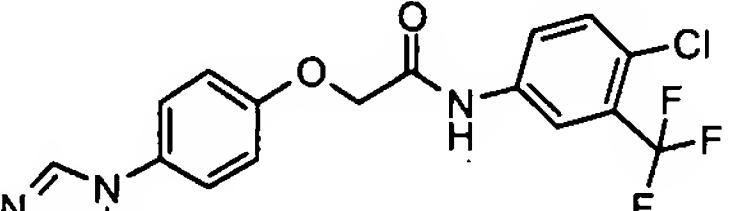
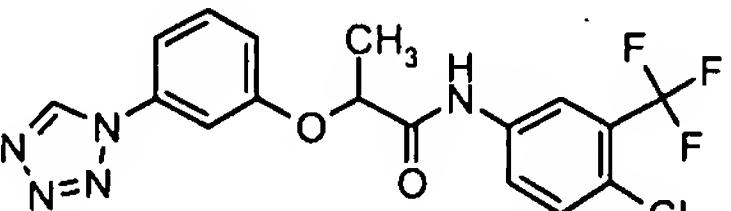
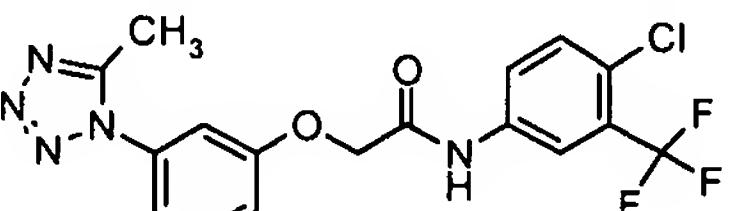
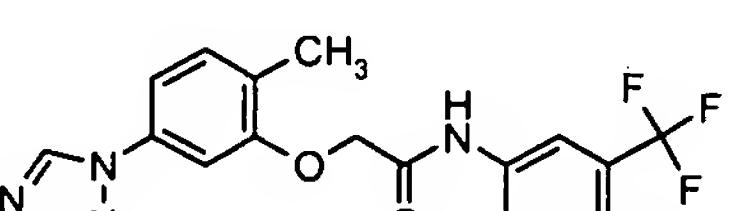
Entry	Name	Structure
14	N-[4-bromo-3-(trifluoromethyl)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
15	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[4-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
16	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}propanamide	
17	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[3-(5-methyl-1H-tetrazol-1-yl)phenyl]oxy}acetamide	
18	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[2-methyl-5-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	

Table 3

Entry	Name	Structure
19	N-(4-chlorophenyl)-N-methyl-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}acetamide	
20	N-[4-chloro-2-(trifluoromethyl)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}acetamide	
21	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{{3-(2,5-dioxopyrrolidin-1-yl)phenyl}oxy}acetamide	
22	(2E)-N-[4-chloro-3-(trifluoromethyl)phenyl]-3-{{3-(1H-tetrazol-1-yl)phenyl}prop-2-enyl}acetamide	
23	N-[4-fluoro-3-(trifluoromethyl)phenyl]-2-{{4-(1H-tetrazol-1-yl)phenyl}oxy}acetamide	
24	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{{3-(2-methyl-2H-tetrazol-5-yl)phenyl}oxy}acetamide	

Table 3

Entry	Name	Structure
25	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[2,4-dichloro-5-(1H-tetrazol-1- yl)phenyl]oxy}acetamide	
26	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[3-(1H-tetrazol-1- yl)phenyl]thio}acetamide	
27	N-[4-chloro-3-(trifluoromethyl)phenyl]- N~2~-[3-(1H-tetrazol-1- yl)phenyl]glycinamide	
28	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[2-(1H-tetrazol-1- yl)phenyl]oxy}acetamide	
29	methyl 1-[(2-[(4-chloro-3-(trifluoromethyl)phenyl)amino]-2- ethoxy)phenyl]-1H-1,2,3-triazole-4- carboxylate	

Table 3

Entry	Name	Structure
30	1,1 dimethylethyl {4 [{[3-(1H tetrazol-1-yl)phenyl]oxy}acetyl]amino}phenyl}carbamate	
31	1,1 dimethylethyl {4 [{[4-(1H tetrazol-1-yl)phenyl]oxy}acetyl]amino}phenyl}carbamate	
32	N-{4-[(1-ethylpiperidin-4-yl)amino]phenyl}-2-[(3-(1H tetrazol-1-yl)phenyl)oxy]acetamide	
33	N-{4-[(1-ethylpiperidin-3-yl)amino]phenyl}-2-[(3-(1H tetrazol-1-yl)phenyl)oxy]acetamide	
34	N-(4-aminophenyl)-2-[(4-(1H tetrazol-1-yl)phenyl)oxy]acetamide	

Table 3

Entry	Name	Structure
35	N-[4-[(1-ethylpiperidin-4-yl)amino]phenyl]-2-[(4-(1H-tetrazol-1-yl)phenyl)oxy]acetamide	
36	N-[4-[(1-ethylpiperidin-3-yl)amino]phenyl]-2-[(4-(1H-tetrazol-1-yl)phenyl)oxy]acetamide	
37	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(3-pyridin-4-ylphenyl)oxy]acetamide	
38	N-[4-chloro-3-(trifluoromethyl)phenyl]-N~2~-methyl-N~2~-[3-(1H-tetrazol-1-yl)phenyl]glycinamide	
39	N-[1,3-benzothiazol-2-yl]-2-[(3-(1H-tetrazol-1-yl)phenyl)oxy]acetamide	

Table 3

Entry	Name	Structure
40	N-quinolin-8-yl-2-([3-(1H-tetrazol-1-yl)phenyl]oxy)acetamide	
41	N-(2,3-dihydro-1,4-benzodioxin-6-yl)-2-([3-(1H-tetrazol-1-yl)phenyl]oxy)acetamide	
42	N-isoquinolin-5-yl-2-([3-(1H-tetrazol-1-yl)phenyl]oxy)acetamide	
43	N-[3-[(phenylmethyl)oxy]phenyl]-2-([3-(1H-tetrazol-1-yl)phenyl]oxy)acetamide	
44	N-[5-methyl-2-(methoxy)phenyl]-2-([3-(1H-tetrazol-1-yl)phenyl]oxy)acetamide	

Table 3

Entry	Name	Structure
45	N-[2,5-bis(methyloxy)phenyl]-2-([3-(1H-tetrazol-1-yl)phenyl]oxy)acetamide	
46	N-(6-fluoro-1,3-benzothiazol-2-yl)-2-([3-(1H-tetrazol-1-yl)phenyl]oxy)acetamide	
47	methyl 3-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetyl]amino]benzoate	
48	5-chloro-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetyl]amino]benzamide	
49	N-[5-chloro-2,4-bis(methyloxy)phenyl]-2-([3-(1H-tetrazol-1-yl)phenyl]oxy)acetamide	

Table 3

Entry	Name	Structure
50	$\text{N-[2-(phenyloxy)phenyl]-2-}\{\text{[3-(1H-tetrazol-1-yl)phenyl]oxy}\}\text{acetamide}$	
51	$\text{N-[3-(aminosulfonyl)phenyl]-2-}\{\text{[3-(1H-tetrazol-1-yl)phenyl]oxy}\}\text{acetamide}$	
52	$\text{N-[2-(methyloxy)-5-(trifluoromethyl)phenyl]-2-}\{\text{[3-(1H-tetrazol-1-yl)phenyl]oxy}\}\text{acetamide}$	
53	$\text{N-(4-}\{\text{[(4-methylphenyl)sulfonyl]amino}\}\text{phenyl)-2-}\{\text{[3-(1H-tetrazol-1-yl)phenyl]oxy}\}\text{acetamide}$	
54	$\text{N-(5-phenyl-1H-pyrazol-3-yl)-2-}\{\text{[3-(1H-tetrazol-1-yl)phenyl]oxy}\}\text{acetamide}$	

Table 3

Entry	Name	Structure
55	N-1,3-benzothiazol-2-yl-2-[(4-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
56	N-quinolin-8-yl-2-[(4-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
57	1,1-dimethylethyl 2-{3-[(2-[(4-chloro-3-(trifluoromethyl)phenyl]amino)-2-oxoethyl]oxy]phenyl}-1H-pyrrole-1-carboxylate	
58	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(3-(1H-pyrrol-2-yl)phenyl]oxy}acetamide	
59	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(3-pyrimidin-5-ylphenyl)oxy]acetamide	

Table 3

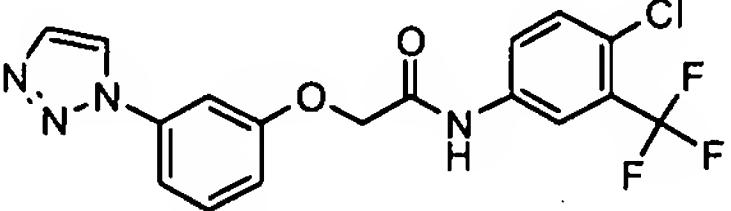
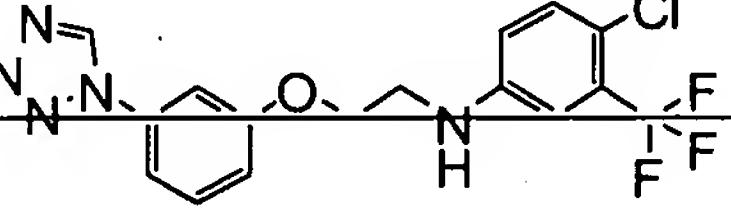
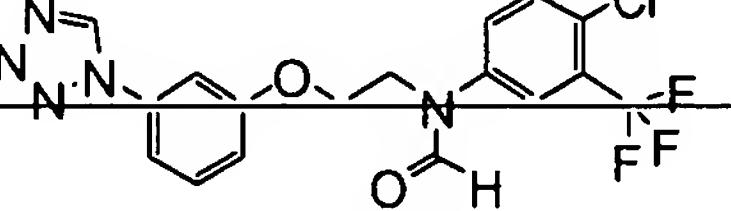
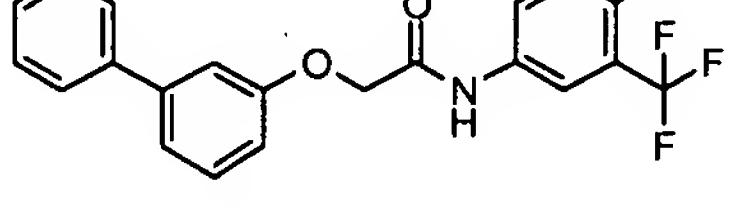
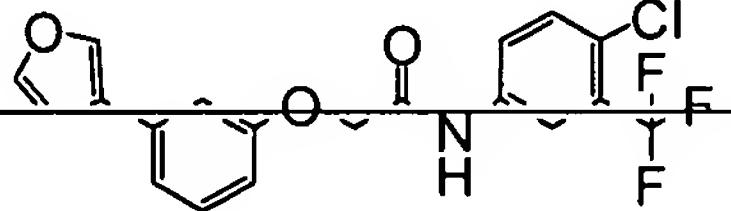
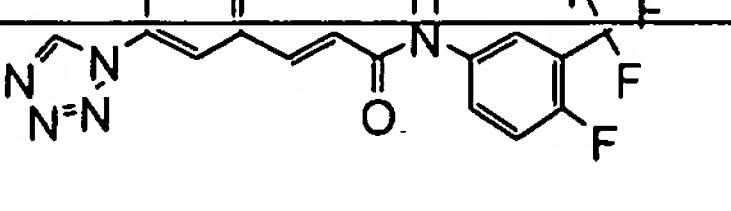
Entry	Name	Structure
60	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[3-(1H-1,2,3-triazol-1- yl)phenyl]oxy}acetamide	
61	4-chloro N-(2-[[3-(1H-tetrazol-1- yl)phenyl]oxy]ethyl)-3- (trifluoromethyl)aniline	
62	N-[4-chloro-3-(trifluoromethyl)phenyl]-N- (2-[[3-(1H-tetrazol-1- yl)phenyl]oxy]ethyl)formamide	
63	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- [(3-pyridin-3-ylphenyl)oxy]acetamide	
64	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- [(3-furan-3-ylphenyl)oxy]acetamide	
65	(2E)-N-[4-fluoro-3- (trifluoromethyl)phenyl]-3-[3-(1H- tetrazol-1-yl)phenyl]prop-2-enamide	

Table 3

Entry	Name	Structure
66	N-[4-fluoro-3-(trifluoromethyl)phenyl]-3-[3-(1H-tetrazol-1-yl)phenyl]propanamide	
67	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(6-(1H-tetrazol-1-yl)pyrimidin-4-yl)oxy]acetamide	
68	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(3-(3,5-dimethylisoxazol-4-yl)phenyl)oxy]acetamide	
69	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(3-quinolin-7-ylphenyl)oxy]acetamide	
70	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(3-furan-2-ylphenyl)oxy]acetamide	

Table 3

Entry	Name	Structure
71	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[3-(1H-tetrazol-1-yl)phenyl]hydrazinecarboxamide	
72	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[3-(dibenzo[b,d]furan-4-ylphenyl)oxy]acetamide	
73	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(4-pyrimidin-5-ylphenyl)oxy]acetamide	
74	N-methyl N-[4-(methoxy)phenyl]-2-[(3-(1H-tetrazol-1-yl)phenyl)oxy]acetamide	
75	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[3-(1H-tetrazol-1-yl)phenyl]methyl}urea	

Table 3

Entry	Name	Structure
76	N-[4-chloro-3-(trifluoromethyl)phenyl]-N-methyl-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide	
77	N-[4-fluoro-3-(trifluoromethyl)phenyl]-N-2-[[3-(1H-tetrazol-1-yl)phenyl]glycinamide	
78	N-[4-fluoro-3-(trifluoromethyl)phenyl]-2-[[3-(pyridin-2-ylamino)phenyl]oxy]acetamide	
79	N-[2-fluoro-5-(trifluoromethyl)phenyl]-2-[[3-(1H-tetrazol-1-yl)phenyl]hydrazinecarboxamide	
80	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(4-pyridin-3-ylphenyl)oxy]acetamide	

Table 3

Entry	Name	Structure
81	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-(3-pyrimidin-5-ylphenyl)methyl]urea	
82	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-(4-pyrimidin-5-ylphenyl)methyl]urea	
83	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-(4-pyridin-3-ylphenyl)methyl]urea	
84	[3-(1H-tetrazol-1-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
85	N-[4-fluoro-3-(trifluoromethyl)phenyl]-2-[(4-pyrimidin-5-ylphenyl)oxy]acetamide	

Table 3

Entry	Name	Structure
86	N~2~-[4-chloro-3-(trifluoromethyl)phenyl]-N-[3-(1H-tetrazol-1-yl)phenyl]glycinamide	
87	2-{{4-chloro-3-(trifluoromethyl)phenyl}oxy}-N-[3-(1H-tetrazol-1-yl)phenyl]acetamide	
88	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{{[3-methyl-4-(1H-tetrazol-1-yl)phenyl]oxy}acetamide}	
89	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{{[4-(1H-1,2,3-triazol-1-yl)phenyl]oxy}acetamide}	
90	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{{[3-fluoro-4-(1H-tetrazol-1-yl)phenyl]oxy}acetamide}	

Table 3

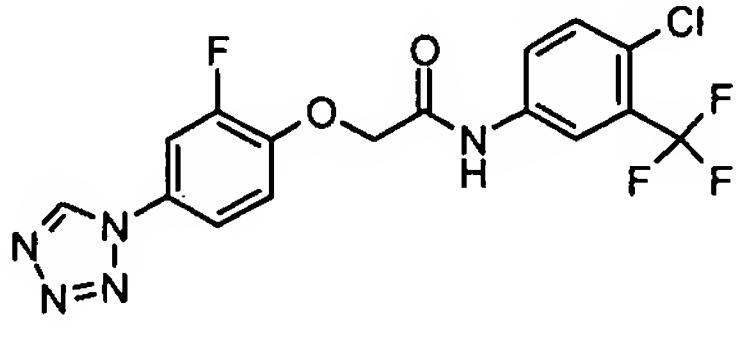
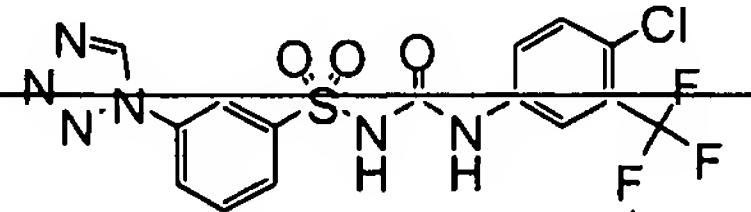
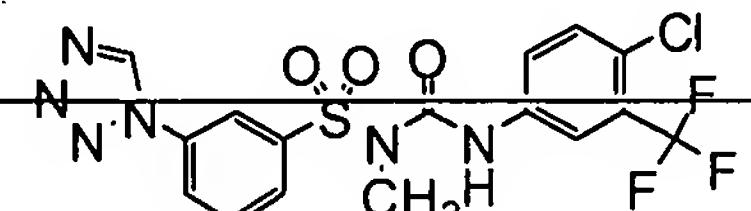
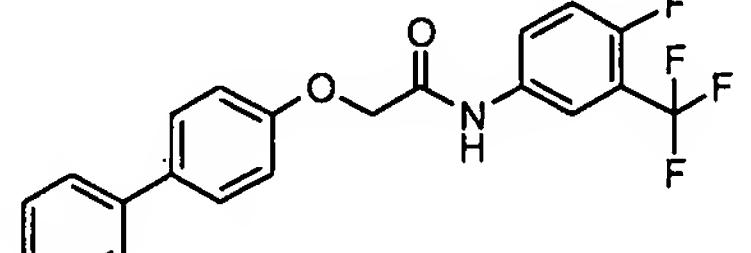
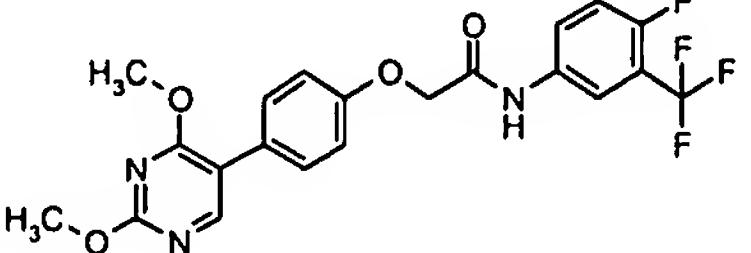
Entry	Name	Structure
91	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[2-fluoro-4-(1H-tetrazol-1- yl)phenyl]oxy}acetamide	
92	N-({[4-chloro-3-(trifluoromethyl)phenyl]amino}carbonyl)-3-(1H-tetrazol-1-yl)benzenesulfonamide	
93	N-({[4-chloro-3-(trifluoromethyl)phenyl]amino}carbonyl)-N-methyl-3-(1H-tetrazol-1-yl)benzenesulfonamide	
94	N-[4-fluoro-3-(trifluoromethyl)phenyl]-2- [(4-pyridin-3-ylphenyl)oxy]acetamide	
95	2-({4-[2,4-bis(methoxy)pyrimidin-5- yl]phenyl}oxy)-N-[4-fluoro-3-(trifluoromethyl)phenyl]acetamide	

Table 3

Entry	Name	Structure
96	2-({4-[2,4-bis(methyloxy)pyrimidin-5-yl]phenyl}oxy)-N-[4-chloro-3-(trifluoromethyl)phenyl]acetamide	
97	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(4-pyridin-4-ylphenyl)oxy]acetamide	
98	N-[4-chloro-3-(trifluoromethyl)phenyl]-N~2~-[3-(methyloxy)-4-(1H-tetrazol-1-yl)phenyl]glycinamide	
99	N-[4-chloro-3-(trifluoromethyl)phenyl]-N~2~-[4-(methyloxy)-3-(1H-tetrazol-1-yl)phenyl]glycinamide	
100	N-[4-chloro-3-(trifluoromethyl)phenyl]-N~2~-[4-(1H-tetrazol-1-yl)phenyl]glycinamide	

Table 3

Entry	Name	Structure
101	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(2,3,5,6-tetrafluoro-4-pyrimidin-5-ylphenyl)hydrazinecarboxamide	
102	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-(4-(1H-tetrazol-1-yl)phenyl)methylurea	
103	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(4-pyrimidin-5-ylphenyl)hydrazinecarboxamide	
104	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-(3-pyridin-3-ylphenyl)methylurea	
105	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-methyl-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}propanamide	

Table 3

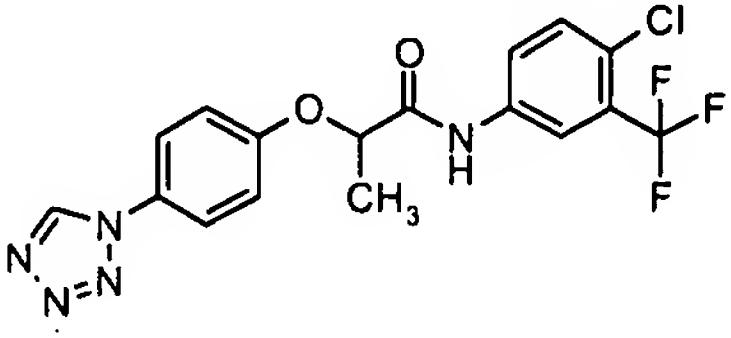
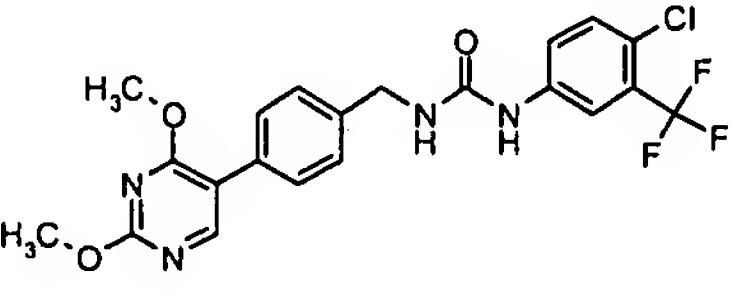
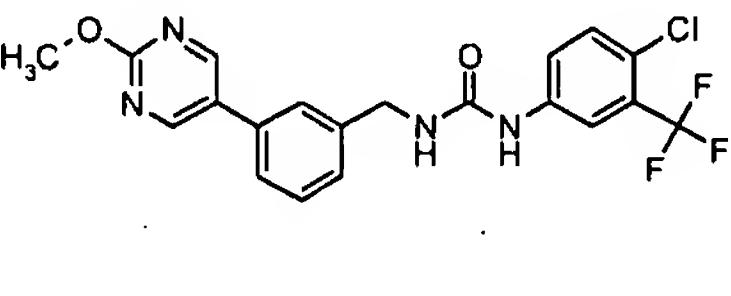
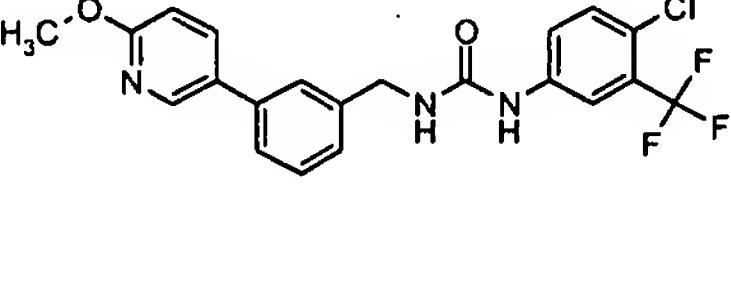
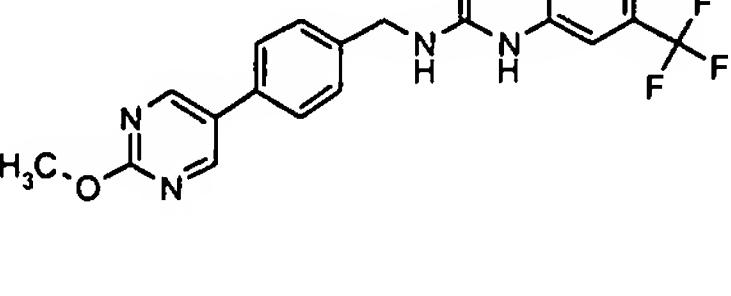
Entry	Name	Structure
106	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[4-(1H-tetrazol-1- yl)phenyl]oxy}propanamide	
107	N-((4-[2,4-bis(methyloxy)pyrimidin-5- yl]phenyl)methyl)-N'-[4-chloro-3- (trifluoromethyl)phenyl]urea	
108	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'- ({3-[2-(methyloxy)pyrimidin-5- yl]phenyl}methyl)urea	
109	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'- ({3-[6-(methyloxy)pyridin-3- yl]phenyl}methyl)urea	
110	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'- ({4-[2-(methyloxy)pyrimidin-5- yl]phenyl}methyl)urea	

Table 3

Entry	Name	Structure
111	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-(4-[6-(methyloxy)pyridin-3-yl]phenyl)methylurea	
112	1,1-dimethylethyl 2-[4-[(2-[(4-chloro-3-(trifluoromethyl)phenyl)amino]-2-oxoethyl)oxy]phenyl]-1H-indole-1-carboxylate	
113	N-[(4-chloro-3-(trifluoromethyl)phenyl)amino]carbonyl)-4-(1H-tetrazol-1-yl)benzenesulfonamide	
114	N-[4-chloro-3-(trifluoromethyl)phenyl]-N-[3-(2H-tetrazol-5-yl)phenyl]glycinamide	
115	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[[2,6-difluoro-4-(1H-tetrazol-1-yl)phenyl]oxy]acetamide	

Table 3

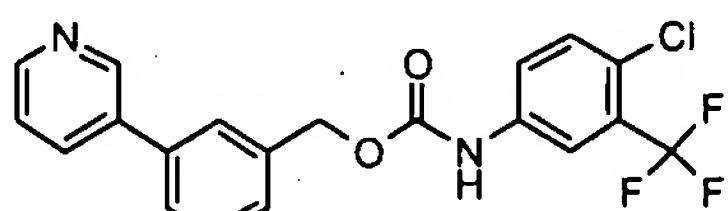
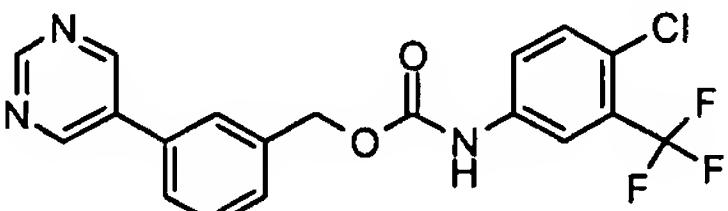
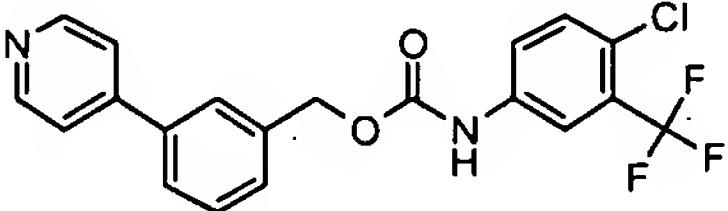
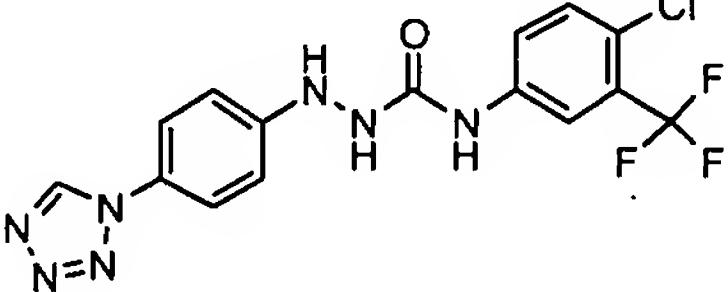
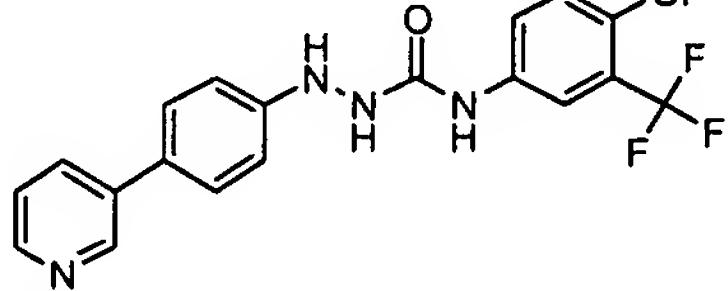
Entry	Name	Structure
116	(3-pyridin-3-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
117	(3-pyrimidin-5-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
118	(3-pyridin-4-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
119	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[4-(1H-tetrazol-1-yl)phenyl]hydrazinecarboxamide	
120	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(4-pyridin-3-ylphenyl)hydrazinecarboxamide	

Table 3

Entry	Name	Structure
121	(4-pyridin-3-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
122	(4-pyridin-4-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
123	(4-pyrimidin-5-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
124	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-(4-pyridin-4-ylphenyl)methylurea	
125	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(3-pyridin-3-ylphenyl)hydrazinecarboxamide	

Table 3

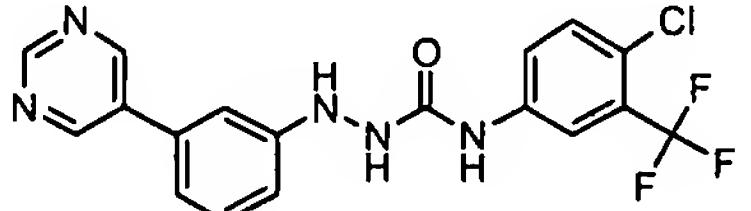
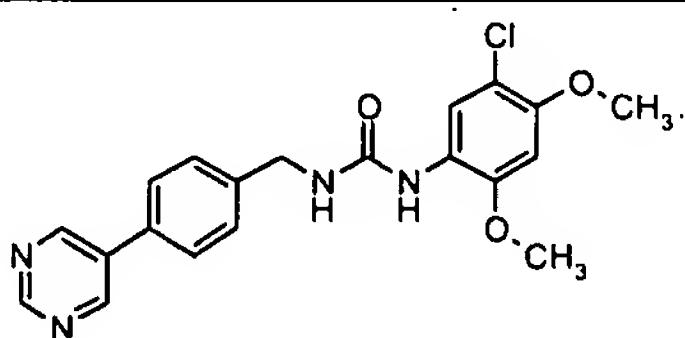
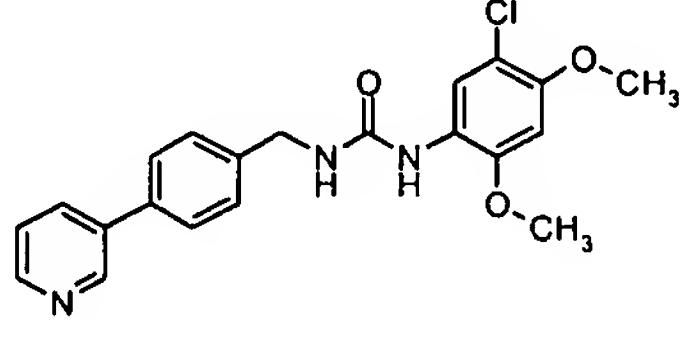
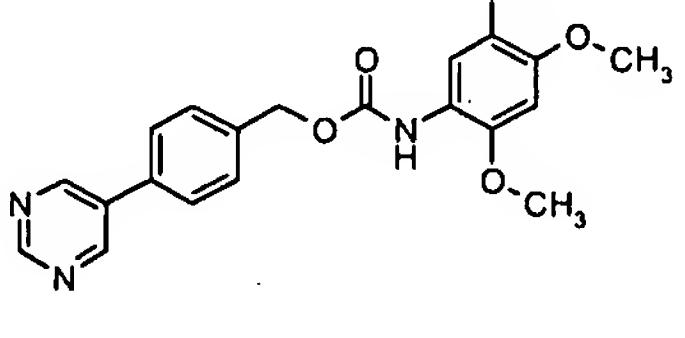
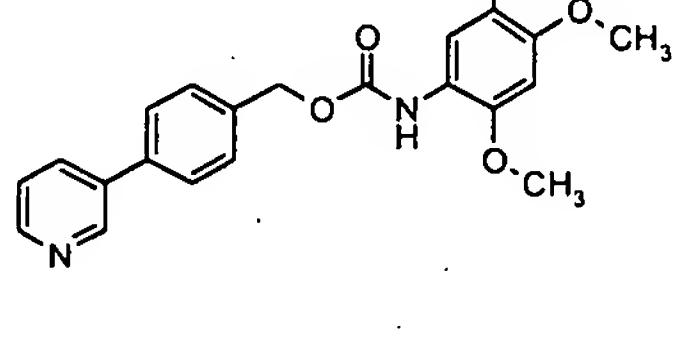
Entry	Name	Structure
126	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(3-pyrimidin-5-ylphenyl)hydrazinecarboxamide	
127	N-[5-chloro-2,4-bis(methyloxy)phenyl]-N'-(4-pyrimidin-5-ylphenyl)methylurea	
128	N-[5-chloro-2,4-bis(methyloxy)phenyl]-N'-(4-pyridin-3-ylphenyl)methylurea	
129	(4-pyrimidin-5-ylphenyl)methyl [5-chloro-2,4-bis(methyloxy)phenyl]carbamate	
130	(4-pyridin-3-ylphenyl)methyl [5-chloro-2,4-bis(methyloxy)phenyl]carbamate	

Table 3

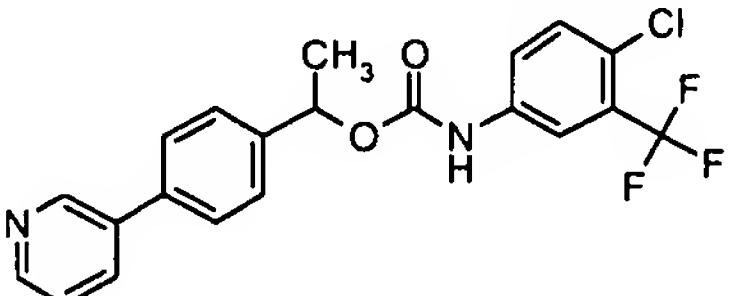
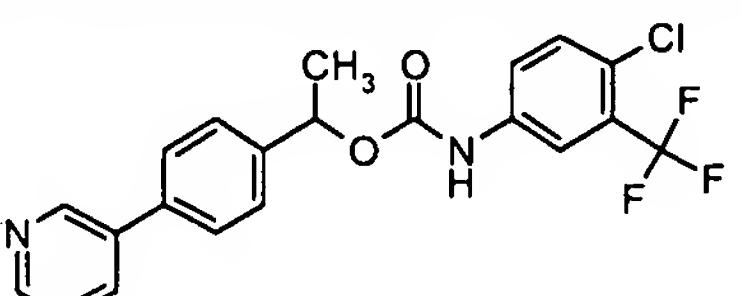
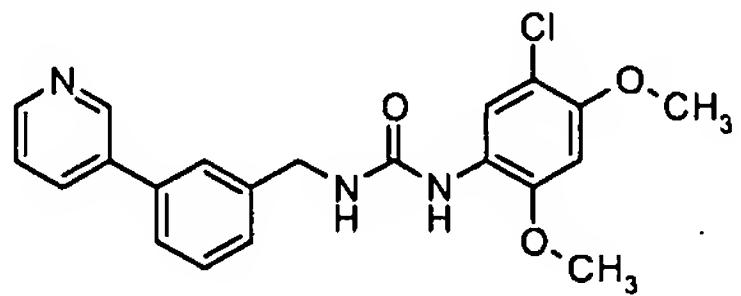
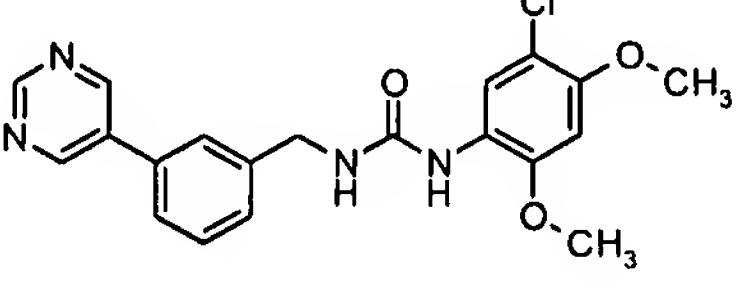
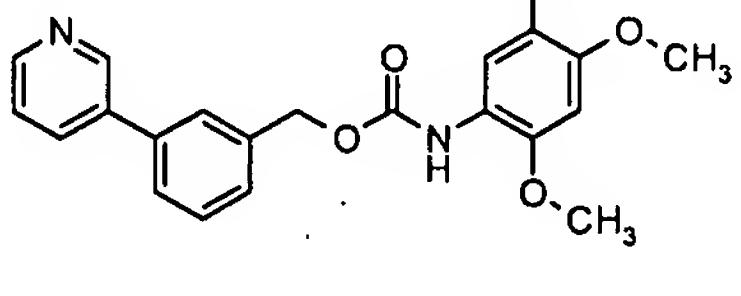
Entry	Name	Structure
131	1-(4-pyridin-3-ylphenyl)ethyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
132	1-(4-pyrimidin-5-ylphenyl)ethyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
133	N-[5-chloro-2,4-bis(methyloxy)phenyl]-N'-(3-pyridin-3-ylphenyl)methylurea	
134	N-[5-chloro-2,4-bis(methyloxy)phenyl]-N'-(3-pyrimidin-5-ylphenyl)methylurea	
135	(3-pyridin-3-ylphenyl)methyl [5-chloro-2,4-bis(methyloxy)phenyl]carbamate	

Table 3

Entry	Name	Structure
136	(3-pyrimidin-5-ylphenyl)methyl [5-chloro-2,4-bis(methyloxy)phenyl]carbamate	
137	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-methyl-2-(3-pyrimidin-5-ylphenyl)hydrazinecarboxamide	
138	N-[4-fluoro-3-(trifluoromethyl)phenyl]-N'-(4-pyridin-3-ylphenyl)methylurea	
139	N-[3-(6-aminopyridin-3-yl)phenyl]methyl-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
140	N-[4-(6-aminopyridin-3-yl)phenyl]methyl-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
141	N-[3-(2-aminopyrimidin-5-yl)phenyl]methyl-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	

Table 3

Entry	Name	Structure
142	$\text{N}-(\text{4-(2-aminopyrimidin-5-yl)phenylmethyl})\text{N}'-(\text{4-chloro-3-(trifluoromethyl)phenyl})\text{urea}$	
143	$\text{N}-(\text{4-chloro-3-(trifluoromethyl)phenyl})\text{-N}'-(\text{1-(4-pyridin-3-ylphenyl)ethyl})\text{urea}$	
144	$\text{N}-(\text{4-chloro-3-(trifluoromethyl)phenyl})\text{-N}'-(\text{1-(4-pyrimidin-5-ylphenyl)ethyl})\text{urea}$	
145	$\text{N}-(\text{4-chloro-3-(trifluoromethyl)phenyl})\text{-2-}(\text{[4-(1H-indol-2-yl)phenyl]oxy})\text{acetamide}$	
146	$\text{N}-(\text{4-chloro-3-(trifluoromethyl)phenyl})\text{-2-}(\text{[isoquinolin-7-yloxy})\text{acetamide}$	

Table 3

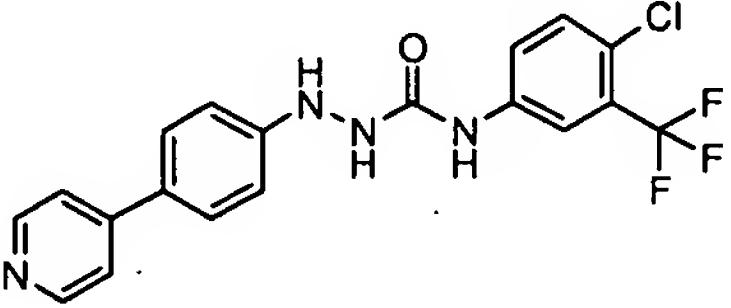
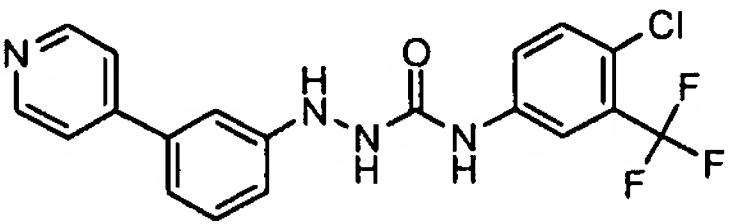
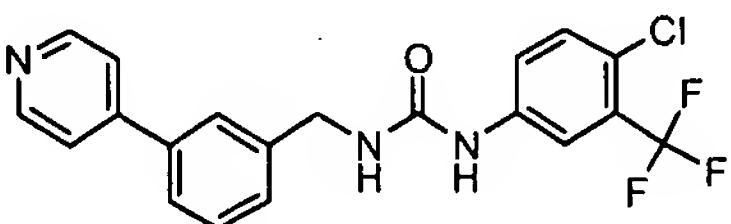
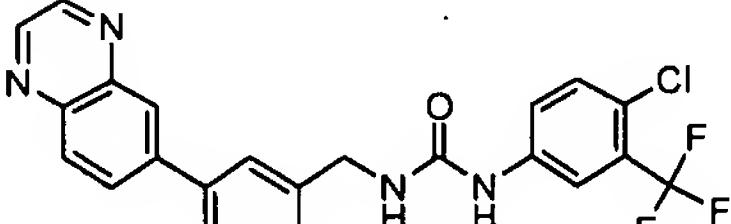
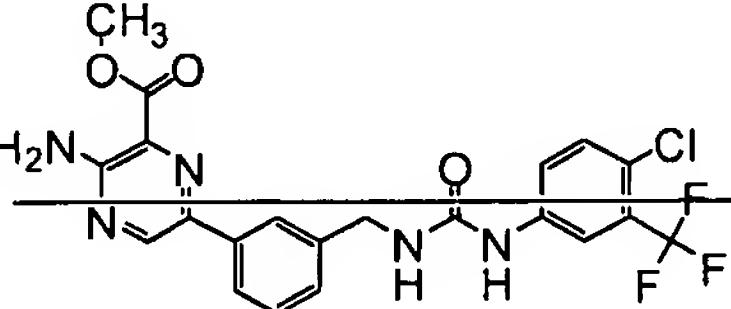
Entry	Name	Structure
147	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(4-pyridin-4-ylphenyl)hydrazinecarboxamide	
148	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(3-pyridin-4-ylphenyl)hydrazinecarboxamide	
149	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'[(3-pyridin-4-ylphenyl)methyl]urea	
150	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'[(3-quinoxalin-6-ylphenyl)methyl]urea	
151	methyl 3-amino-6-((4-chloro-3-(trifluoromethyl)phenyl)amino)carbonyl)amino)methylphenyl)pyrazine-2-carboxylate	

Table 3

Entry	Name	Structure
152	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-(4-quinoxalin-6-ylphenyl)methyl]urea	
153	N-[3-(2-amino-5-methylpyridin-3-yl)phenyl]methyl]-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
154	methyl 3-amino-6-((4-((4-chloro-3-(trifluoromethyl)phenyl)amino)carbonyl)aminomethyl)phenyl)pyrazine-2-carboxylate	
155	[3-(1H-tetrazol-1-yl)phenyl]methyl [3-chloro-4-(methyloxy)phenyl]carbamate	
156	N-[3-chloro-4-(methyloxy)phenyl]-N'-{[3-(1H-tetrazol-1-yl)phenyl]methyl}urea	

Table 3

Entry	Name	Structure
157	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[4-(5-hydroxy-1H-tetrazol-1- yl)phenyl]oxy}acetamide	
158	N-{[3-(2-amino-5-chloropyridin-3- yl)phenyl]methyl}-N'-[4-chloro-3- (trifluoromethyl)phenyl]urea	
159	N-{[4-(2-amino-5-chloropyridin-3- yl)phenyl]methyl}-N'-[4-chloro-3- (trifluoromethyl)phenyl]urea	
160	N-{[3-(6-chloropyridin-3- yl)phenyl]methyl}-N'-[4-chloro-3- (trifluoromethyl)phenyl]urea	
161	N-{[4-(6-chloropyridin-3- yl)phenyl]methyl}-N'-[4-chloro-3- (trifluoromethyl)phenyl]urea	
162	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'- {[4-(pyrimidin-2- yloxy)phenyl]methyl}urea	

Table 3

Entry	Name	Structure
163	$\text{N}(\{[4\text{-chloro-3-}(\text{trifluoromethyl})\text{phenyl}]\text{amino}\}\text{carbonyl})\text{-3-(1H-tetrazol-1-yl)benzamide}$	
164	$\text{3-amino-6-}(\{[4\text{-chloro-3-}(\text{trifluoromethyl})\text{phenyl}]\text{amino}\}\text{carbonyl})\text{-2-}(\text{dimethylamino})\text{ethyl}\text{-}N\{2-}(\text{dimethylamino})\text{ethyl}\text{-}N\{2-}\text{carboxamide}$	
165	$\text{N-[4-chloro-3-(trifluoromethyl)phenyl]-N'}\{[3-(6-fluoropyridin-3-yl)phenyl]\text{methyl}\}\text{urea}$	
166	$\text{N-[4-chloro-3-(trifluoromethyl)phenyl]-N'}\{[3-[2-(methyloxy)pyridin-3-yl]phenyl]\text{methyl}\}\text{urea}$	
167	$\text{N-[4-chloro-3-(trifluoromethyl)phenyl]-N'}\{[4-(6-fluoropyridin-3-yl)phenyl]\text{methyl}\}\text{urea}$	

Table 3

Entry	Name	Structure
168	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-(4-[2-(methyloxy)pyridin-3-yl]phenyl)methyl)urea	
169	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-(4-(6-methylpyridin-3-yl)phenyl)methyl)urea	
170	N-[4-(2-amino-5-fluoropyridin-3-yl)phenyl]methyl)-N'-(4-chloro-3-(trifluoromethyl)phenyl)urea	
171	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-(3-(6-methylpyridin-3-yl)phenyl)methyl)urea	
172	N-[4-(2-aminopyridin-3-yl)phenyl]methyl)-N'-(4-chloro-3-(trifluoromethyl)phenyl)urea	
173	N-[3-(2-aminopyridin-3-yl)phenyl]methyl)-N'-(4-chloro-3-(trifluoromethyl)phenyl)urea	

Table 3

Entry	Name	Structure
174	[3-(6-methylpyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
175	[3-(2-amino-5-fluoropyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
176	[3-(2-aminopyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
177	(3-pyrazin-2-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
178	N-[4-chloro-3-(trifluoromethyl)phenyl]N'-(3-[6-(hydroxymethyl)pyridin-3-yl]phenyl)methylurea	
179	N-{{[3-(6-acetylpyridin-3-yl)phenyl]methyl}-N'-(4-chloro-3-(trifluoromethyl)phenyl)urea	

Table 3

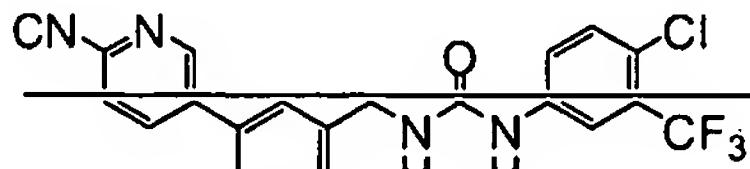
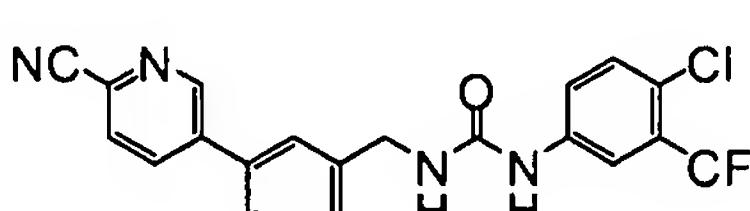
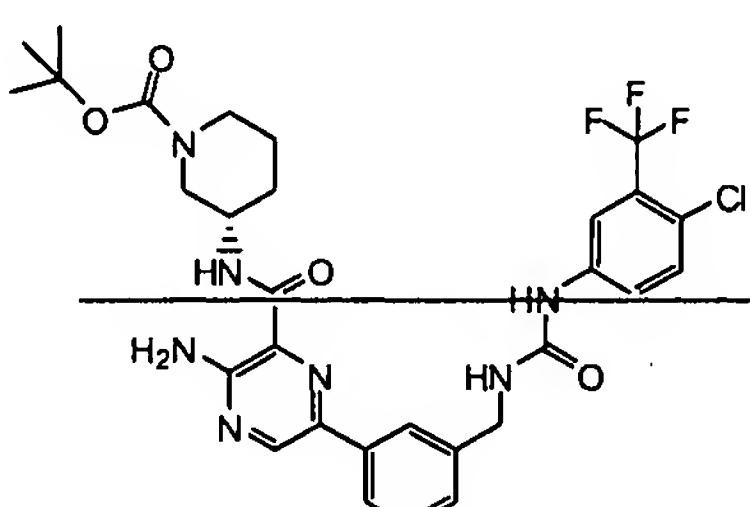
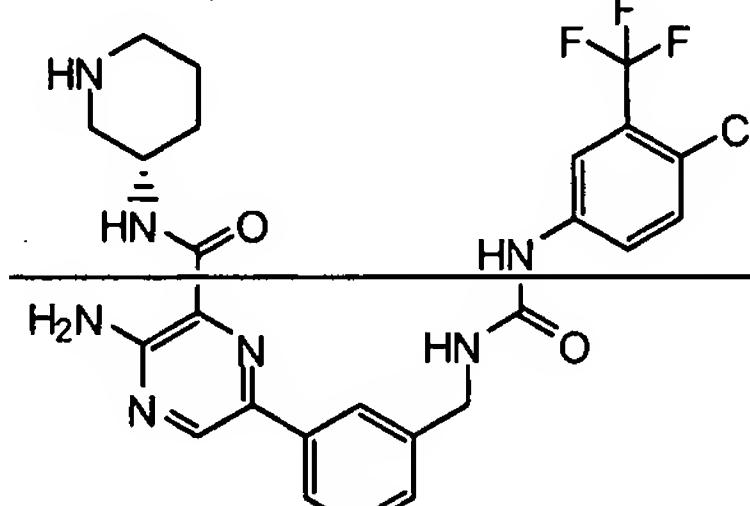
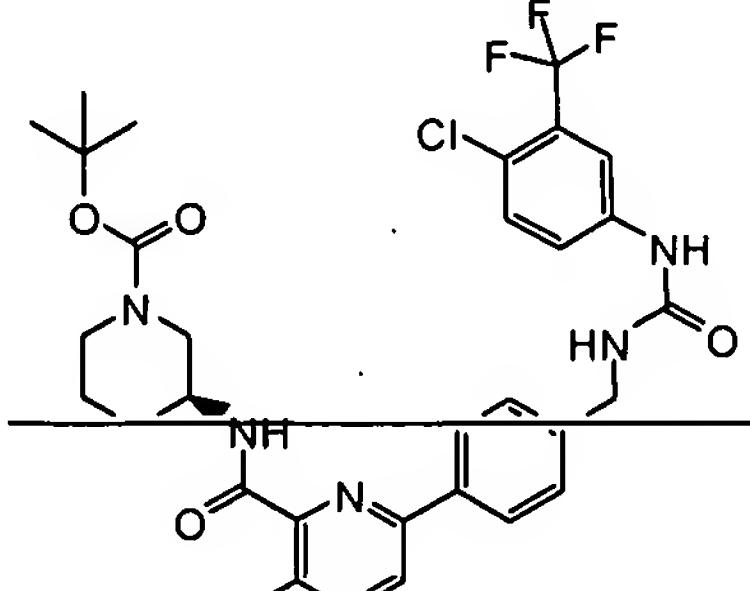
Entry	Name	Structure
180	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[3-(6-cyanopyridin-3-yl)phenyl]methyl}urea	 
181	1,1-dimethylethyl (3S)-3-({{3-amino-6-(3-((4-chloro-3-(trifluoromethyl)phenyl)amino)carbonyl)amino}methyl}phenyl)pyrazin-2-ylcarbonyl)amino)piperidine-1-carboxylate	
182	3-amino-6-(3-((4-chloro-3-(trifluoromethyl)phenyl)amino)carbonyl)amino)methyl)phenyl)N-[(3S)-piperidin-3-yl]pyrazine-2-carboxamide	
183	1,1-dimethylethyl (3S)-3-({{3-amino-6-(4-(4-chloro-3-(trifluoromethyl)phenyl)amino)carbonyl)amino}methyl}phenyl)pyrazin-2-ylcarbonyl)amino)piperidine-1-carboxylate	

Table 3

Entry	Name	Structure
184	3-amino-6-(4-((4-chloro-3-(trifluoromethyl)phenyl)amino)carbonyl)amino]methyl]phenyl)N-(3S)-piperidin-3-yl]pyrazine-2-carboxamide	
185	[3-(7H-pyrrolo[2,3-d]pyrimidin-4-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
186	N-((3-(2-amino-5-fluoropyridin-3-yl)phenyl)methyl)-N'-(4-chloro-3-(trifluoromethyl)phenyl)urea	
187	[6-(1H-tetrazol-1-yl)pyridin-2-yl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
188	[3-(1H-benzimidazol-2-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	

Table 3

Entry	Name	Structure
189	[3-(6-amino-2-methylpyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
190	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-(3-[5-(methylthio)pyridin-3-yl]phenyl)methylurea	
191	[4-(6-methylpyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
192	[4-(2-amino-5-fluoropyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
193	[4-(2-aminopyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
194	(4-pyrazin-2-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	

Table 3

Entry	Name	Structure
195	[4-(7H-pyrrolo[2,3-d]pyrimidin-4-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
196	[4-(6-amino-2-methylpyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
197	[3-(1H-tetrazol-1-yl)phenyl]methyl 1,3-benzothiazol-2-ylcarbamate	
198	[3-(1H-tetrazol-1-yl)phenyl]methyl (5-bromopyridin-2-yl)carbamate	
199	(3-pyridin-3-ylphenyl)methyl (3,5-dimethylphenyl)carbamate	
200	(3-pyridin-3-ylphenyl)methyl [5-chloro-2-(methyloxy)phenyl]carbamate	

Table 3

Entry	Name	Structure
201	[4-(1H-tetrazol-1-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
202	(3-pyrimidin-5-ylphenyl)methyl [5-chloro-2-(methyloxy)phenyl]carbamate	
203	(4-pyrimidin-5-ylphenyl)methyl (3,4-dimethylphenyl)carbamate	
204	(3-pyridin-3-ylphenyl)methyl (3,4-dimethylphenyl)carbamate	
205	1,1-dimethyl-3-((3-amino-6-((4-chloro-3-(trifluoromethyl)phenyl)amino)carbonyl)amino)methylphenyl)pyrazin-2-ylcarbonyl)amino)piperidine-1-carboxylate	

Table 3

Entry	Name	Structure
206	1,1-dimethylethyl 3-((3-amino-6-(4-((4-chloro-3-(trifluoromethyl)phenyl)amino)carbonyl)amino)methyl)phenyl)pyrazin-2-ylcarbonyl)amino)piperidine-1-carboxylate	
207	3-amino-6-(3-((4-chloro-3-(trifluoromethyl)phenyl)amino)carbonyl)amino)methyl)phenyl)N-piperidin-3-ylpyrazine-2-carboxamide	
208	3-amino-6-(4-((4-chloro-3-(trifluoromethyl)phenyl)amino)carbonyl)amino)methyl)phenyl)N-piperidin-3-ylpyrazine-2-carboxamide	

Table 3

Entry	Name	Structure
209	1,1-dimethylethyl 4-[[3-amino-6-(3-((4-chloro-3-(trifluoromethyl)phenyl)amino)carbonyl)amino]methyl]phenyl]pyrazin-2-yl]carbonyl)piperazine-1-carboxylate	
210	1,1-dimethylethyl 4-[[3-amino-6-(4-((4-chloro-3-(trifluoromethyl)phenyl)amino)carbonyl)amino]methyl]phenyl]pyrazin-2-yl]carbonyl)piperazine-1-carboxylate	
211	N-((3-[5-amino-6-(piperazin-1-yl)carbonyl]pyrazin-2-yl)phenyl)methyl)-N'-(4-chloro-3-(trifluoromethyl)phenyl)urea	

Table 3

Entry	Name	Structure
212	<p>N-({4-[5-amino-6-(piperazin-1-yl)carbonyl]pyrazin-2-yl}phenyl)methyl)-</p> <p>N'-[4-chloro-3-(trifluoromethyl)phenyl]urea</p>	
213	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[3-(1H-pyrazol-4-yl)phenyl]methyl}urea	
214	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[4-(1H-pyrazol-4-yl)phenyl]methyl}urea	
215	[3-(2-piperazin-1-ylpyrimidin-5-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	

Table 3

Entry	Name	Structure
216	[4-(2-piperazin-1-ylpyrimidin-5-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
217	N- {[3-(2-chloropyridin-3-yl)phenyl]methyl} -N' -[4-chloro-3-(trifluoromethyl)phenyl]urea	
218	N- {[4-(2-chloropyridin-3-yl)phenyl]methyl} -N' -[4-chloro-3-(trifluoromethyl)phenyl]urea	
219	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[3-(2-fluoropyridin-3-yl)phenyl]methyl}urea	
220	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[4-(2-fluoropyridin-3-yl)phenyl]methyl}urea	

Table 3

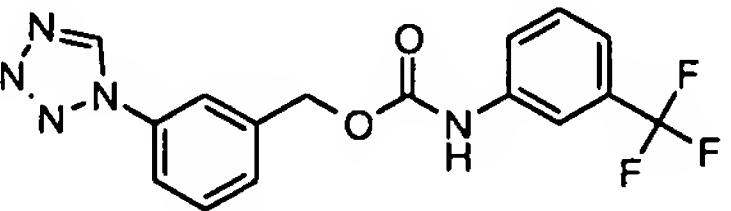
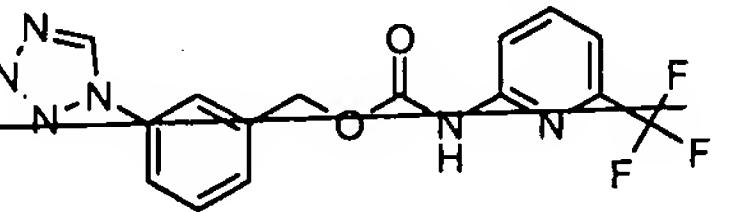
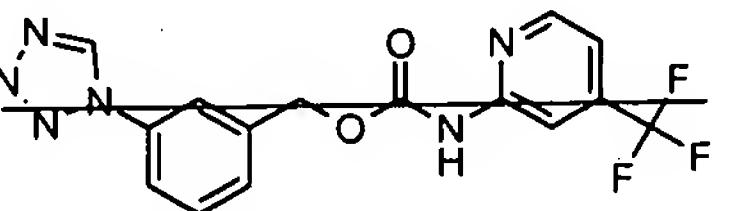
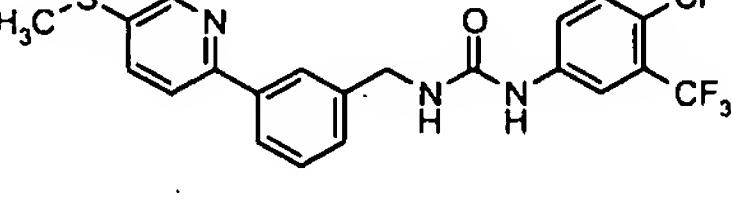
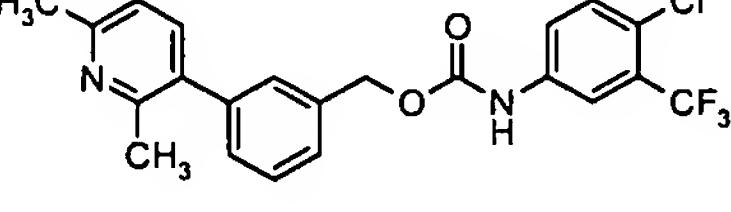
Entry	Name	Structure
221	[3-(1H-tetrazol-1-yl)phenyl]methyl [3-(trifluoromethyl)phenyl]carbamate	
222	[3-(1H-tetrazol-1-yl)phenyl]methyl [6-(trifluoromethyl)pyridin-2-yl]carbamate	
223	[3-(1H-tetrazol-1-yl)phenyl]methyl [4-(trifluoromethyl)pyridin-2-yl]carbamate	
224	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-(3-[5-(methylthio)pyridin-2-yl]phenyl)methylurea	
225	[3-(2,6-dimethylpyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	

Table 3

Entry	Name	Structure
226	{3-[5-(methyloxy)pyridin-3-yl]phenyl}methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
227	2,3'-bipyridin-6-ylmethyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
228	(6-pyrimidin-5-ylpyridin-2-yl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
229	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-(3-isoquinolin-4-ylphenyl)methyl]urea	
230	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-(4-isoquinolin-4-ylphenyl)methyl]urea	

Table 3

Entry	Name	Structure
231	[6-(1H-tetrazol-1-yl)pyridin-2-yl]methyl [4-(trifluoromethyl)pyridin-2-yl]carbamate	
232	[3-(1H-pyrazol-4-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
233	[4-(1H-pyrazol-4-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	

28. (previously presented) A pharmaceutical composition comprising the compound according to claim 1 and a pharmaceutically acceptable carrier.
29. (cancelled)
30. (withdrawn from consideration, currently amended) A method for modulating the *in-vivo* activity of a kinase, the method comprising administering to a subject an effective amount of the compound according to claim 1. ~~claim 1 or a compound selected from N-naphthalen-1-yl-2-([3-(1H-tetrazol-1-yl)phenyl]oxy) acetamide, N-[4(phenyloxy)phenyl]-2-([3-(1H-tetrazol-1-yl)phenyl]oxy) acetamide, N-(3,4-dimethylphenyl)-2-([3-(1H-tetrazol-1-yl)phenyl]oxy)~~

acetamide, $\text{N}(\text{2,3-dimethylphenyl})2\{[\text{3-(1H-tetrazol-1-yl)phenyl}]\text{oxy}\}$ acetamide, $\text{N}(\text{2,4-dimethylphenyl})2\{[\text{3-(1H-tetrazol-1-yl)phenyl}]\text{oxy}\}$ acetamide, $\text{N}(\text{2,5-dimethylphenyl})2\{[\text{3-(1H-tetrazol-1-yl)phenyl}]\text{oxy}\}$ acetamide, $\text{N}(\text{3,5-dimethylphenyl})2\{[\text{3-(1H-tetrazol-1-yl)phenyl}]\text{oxy}\}$ acetamide, $\text{N}(\text{2,6-dimethylphenyl})2\{[\text{3-(1H-tetrazol-1-yl)phenyl}]\text{oxy}\}$ acetamide, $\text{N}(\text{2,4,6-trimethylphenyl})\text{acetamide}$, $\text{N}(\text{2-ethylphenyl})2\{[\text{3-(1H-tetrazol-1-yl)phenyl}]\text{oxy}\}$ acetamide, $\text{N}(\text{4-ethylphenyl})2\{[\text{3-(1H-tetrazol-1-yl)phenyl}]\text{oxy}\}$ acetamide, $\text{N}(\text{2,6-diethylphenyl})2\{[\text{3-(1H-tetrazol-1-yl)phenyl}]\text{oxy}\}$ acetamide, $\text{N}[\text{2-(methyloxy)phenyl}]2\{[\text{3-(1H-tetrazol-1-yl)phenyl}]\text{oxy}\}$ acetamide, $\text{N}[\text{2-(ethyloxy)phenyl}]2\{[\text{3-(1H-tetrazol-1-yl)phenyl}]\text{oxy}\}$ acetamide, $\text{N}[\text{3-(ethyloxy)phenyl}]2\{[\text{3-(1H-tetrazol-1-yl)phenyl}]\text{oxy}\}$ acetamide, $\text{N}[\text{2,4-bis(methyloxy)phenyl}]2\{[\text{3-(1H-tetrazol-1-yl)phenyl}]\text{oxy}\}$ acetamide, $\text{N}[\text{4-(dimethylamino)phenyl}]2\{[\text{3-(1H-tetrazol-1-yl)phenyl}]\text{oxy}\}$ acetamide, $\text{N}(\text{2,3-dichlorophenyl})2\{[\text{3-(1H-tetrazol-1-yl)phenyl}]\text{oxy}\}$ acetamide, $\text{N}(\text{4-chloro-3-methylphenyl})2\{[\text{3-(1H-tetrazol-1-yl)phenyl}]\text{oxy}\}$ acetamide, $\text{N}(\text{4-bromophenyl})2\{[\text{3-(1H-tetrazol-1-yl)phenyl}]\text{oxy}\}$ acetamide, $\text{N}(\text{2-fluorophenyl})2\{[\text{3-(1H-tetrazol-1-yl)phenyl}]\text{oxy}\}$ acetamide, $\text{N}(\text{4-fluorophenyl})2\{[\text{3-(1H-tetrazol-1-yl)phenyl}]\text{oxy}\}$ acetamide, $\text{N}[\text{2-(trifluoromethyl)phenyl}]2\{[\text{3-(1H-tetrazol-1-yl)phenyl}]\text{oxy}\}$ acetamide, $\text{N}[\text{3-(trifluoromethyl)phenyl}]2\{[\text{3-(1H-tetrazol-1-yl)phenyl}]\text{oxy}\}$ acetamide, $\text{N}[\text{3-(1H-tetrazol-1-yl)phenyl}]\text{oxy}\}$ acetamide, $\text{N}[\text{3-(1H-tetrazol-1-yl)phenyl}]\text{oxy}\}$ acetamide, $\text{N}[\text{2-chloro-5-(trifluoromethyl)phenyl}]2\{[\text{3-(1H-tetrazol-1-yl)phenyl}]\text{oxy}\}$ acetamide, $\text{N}[\text{4-chloro-3-(trifluoromethyl)phenyl}]2\{[\text{3-(1H-tetrazol-1-yl)phenyl}]\text{oxy}\}$ acetamide, $\text{N}(\text{4-chlorophenyl})2\{[\text{3-(1H-tetrazol-1-yl)phenyl}]\text{oxy}\}$ acetamide, $\text{N}(\text{4-aminophenyl})2\{[\text{3-(1H-tetrazol-1-yl)phenyl}]\text{oxy}\}$ acetamide, $\text{N}(\text{4-acetylphenyl})2\{[\text{3-(1H-tetrazol-1-yl)phenyl}]\text{oxy}\}$ acetamide.

31. (withdrawn from consideration) The method according to claim 30, wherein the kinase is c-Kit.

32. (withdrawn from consideration) The method according to claim 31, wherein modulating the *in vivo* activity of c-Kit comprises inhibition of c-Kit.

33. (withdrawn from consideration, currently amended) A method of treating rheumatoid arthritis, graft-host diseases, multiple sclerosis, psoriasis; atherosclerosis, myocardiocircumflex, ischemia, stroke, restenosis; interbowel diseases, osteoarthritis, macular degeneration, or diabetic retinopathy, diseases or disorders associated with uncontrolled, abnormal, and/or unwanted cellular activities, the method comprising administering, to a mammal in need thereof, a therapeutically effective amount of the compound or the pharmaceutical composition as described in claim 1 or a compound, or a pharmaceutical composition comprising said compound, selected from ~~N-naphthalen-1-yl-2-([3-(1H-tetrazol-1-yl)phenyl]oxy) acetamide, N-[4-(phenyloxy)phenyl]-2-([3-(1H-tetrazol-1-yl)phenyl]oxy) acetamide, N-(3,4-dimethylphenyl)-2-([3-(1H-tetrazol-1-yl)phenyl]oxy) acetamide, N-(2,3-dimethylphenyl)-2-([3-(1H-tetrazol-1-yl)phenyl]oxy) acetamide, N-(2,4-dimethylphenyl)-2-([3-(1H-tetrazol-1-yl)phenyl]oxy) acetamide, N-(2,5-dimethylphenyl)-2-([3-(1H-tetrazol-1-yl)phenyl]oxy) acetamide, N-(3,5-dimethylphenyl)-2-([3-(1H-tetrazol-1-yl)phenyl]oxy) acetamide, N-(2,6-dimethylphenyl)-2-([3-(1H-tetrazol-1-yl)phenyl]oxy) acetamide, 2-([3-(1H-tetrazol-1-yl)phenyl]oxy)-N-(2,4,6-trimethylphenyl) acetamide, N-(2-ethylphenyl)-2-([3-(1H-tetrazol-1-yl)phenyl]oxy) acetamide, N-(4-ethylphenyl)-2-([3-(1H-tetrazol-1-yl)phenyl]oxy) acetamide, N-(2,6-diethylphenyl)-2-([3-(1H-tetrazol-1-yl)phenyl]oxy) acetamide, N-[2-(methyloxy)phenyl]-2-([3-(1H-tetrazol-1-yl)phenyl]oxy) acetamide, N-[2-(ethyloxy)phenyl]-2-([3-(1H-tetrazol-1-yl)phenyl]oxy) acetamide, N-[3-(ethyloxy)phenyl]-2-([3-(1H-tetrazol-1-yl)phenyl]oxy) acetamide, N-[2,4-bis(methyloxy)phenyl]-2-([3-(1H-tetrazol-1-yl)phenyl]oxy) acetamide, N-[4-(dimethylamino)phenyl]-2-([3-(1H-tetrazol-1-yl)phenyl]oxy) acetamide, N-(2,3-dichlorophenyl)-2-([3-(1H-tetrazol-1-yl)phenyl]oxy) acetamide, N-(4-chloro-3-methylphenyl)-2-([3-(1H-tetrazol-1-yl)phenyl]oxy) acetamide, N-(4-bromophenyl)-2-([3-(1H-tetrazol-1-yl)phenyl]oxy) acetamide, N-(2-fluorophenyl)-2-([3-(1H-tetrazol-1-yl)phenyl]oxy) acetamide, N-(4-fluorophenyl)-2-([3-(1H-tetrazol-1-yl)phenyl]oxy) acetamide, 2-([3-(1H-tetrazol-1-yl)phenyl]oxy)-N-[2-(trifluoromethyl)phenyl] acetamide, 2-([3-(1H-tetrazol-1-yl)phenyl]oxy)-N-[3-(trifluoromethyl)phenyl] acetamide, methyl-4-([([3-(1H-tetrazol-1-~~

~~y1)phenyl]oxy}acetyl)amino] benzoate, ethyl 4 [{[3 (1H tetrazol-1-yl)phenyl]oxy}acetyl)amino] benzoic acid, N-[3 (methyloxy)phenyl] 2 {[3 (1H tetrazol-1-yl)phenyl]oxy} acetamide, N-[4 (methyloxy)phenyl] 2 {[3 (1H tetrazol-1-yl)phenyl]oxy} acetamide, N-[2 chloro-5-(trifluoromethyl)phenyl] 2 {[3 (1H tetrazol-1-yl)phenyl]oxy} acetamide, N-[4 chloro-3-(trifluoromethyl)phenyl] 2 {[3 (4H 1,2,4 triazol-4-yl)phenyl]oxy} acetamide, N-(4 chlorophenyl) 2 {[3 (1H tetrazol-1-yl)phenyl]oxy} acetamide, N-(4 aminophenyl) 2 {[3 (1H tetrazol-1-yl)phenyl]oxy} acetamide, and N-(4 acetylphenyl) 2 {[3 (1H tetrazol-1-yl)phenyl]oxy} acetamide.~~

34. (withdrawn from consideration, currently amended) A method of screening for modulators of c-Kit, the method comprising combining the compound according to claim 1 or a compound selected from ~~N-naphthalen-1-yl 2 {[3 (1H tetrazol-1-yl)phenyl]oxy} acetamide, N-[4 (phenyloxy)phenyl] 2 {[3 (1H tetrazol-1-yl)phenyl]oxy} acetamide, N-(3,4-dimethylphenyl) 2 {[3 (1H tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,3-dimethylphenyl) 2 {[3 (1H tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,4-dimethylphenyl) 2 {[3 (1H tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,5-dimethylphenyl) 2 {[3 (1H tetrazol-1-yl)phenyl]oxy} acetamide, N-(3,5-dimethylphenyl) 2 {[3 (1H tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,6-dimethylphenyl) 2 {[3 (1H tetrazol-1-yl)phenyl]oxy} acetamide, 2 {[3 (1H tetrazol-1-yl)phenyl]oxy} N-(2,4,6-trimethylphenyl)acetamide, N-(2 ethylphenyl) 2 {[3 (1H tetrazol-1-yl)phenyl]oxy} acetamide, N-(4 ethylphenyl) 2 {[3 (1H tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,6-diethylphenyl) 2 {[3 (1H tetrazol-1-yl)phenyl]oxy} acetamide, N-[2 (methyloxy)phenyl] 2 {[3 (1H tetrazol-1-yl)phenyl]oxy} acetamide, N-[2 (ethyloxy)phenyl] 2 {[3 (1H tetrazol-1-yl)phenyl]oxy} acetamide, N-[3 (ethyloxy)phenyl] 2 {[3 (1H tetrazol-1-yl)phenyl]oxy} acetamide, N-[2,4-bis(methyloxy)phenyl] 2 {[3 (1H tetrazol-1-yl)phenyl]oxy} acetamide, N-[4 (dimethylamino)phenyl] 2 {[3 (1H tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,3-dichlorophenyl) 2 {[3 (1H tetrazol-1-yl)phenyl]oxy} acetamide, N-(4 chloro-3-methylphenyl) 2 {[3 (1H tetrazol-1-yl)phenyl]oxy} acetamide, N-(4 bromophenyl) 2 {[3 (1H tetrazol-1-yl)phenyl]oxy} acetamide, N-(2 fluorophenyl) 2 {[3 (1H tetrazol-1-yl)phenyl]oxy} acetamide, N-(4 fluorophenyl) 2 {[3 (1H tetrazol-1-yl)phenyl]oxy} acetamide, 2 {[3 (1H~~

~~tetrazol-1-yl)phenyl]oxy} N-[2-(trifluoromethyl)phenyl] acetamide, 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy} N-[3-(trifluoromethyl)phenyl] acetamide, methyl 4-[[[[3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl]amino] benzoate, ethyl 4-[[[[3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl]amino] benzoate, 3-[[[[3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl]amino] benzoic acid, N-[3-(methyloxy)phenyl] 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy] acetamide, N-[4-(methyloxy)phenyl] 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy] acetamide, N-[2-chloro-5-(trifluoromethyl)phenyl] 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy] acetamide, N-[4-chloro-3-(trifluoromethyl)phenyl] 2-[[3-(4H-1,2,4-triazol-4-yl)phenyl]oxy] acetamide, N-(4-chlorophenyl) 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy] acetamide, N-(4-aminophenyl) 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy] acetamide, and N-(4-acetylphenyl) 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy] acetamide, and at least one candidate agent and determining the effect of the candidate agent on c-Kit activity.~~

35. (withdrawn from consideration, currently amended) A method of inhibiting proliferative activity in a cell, the method comprising administering an effective amount of a composition comprising the compound according to claim 1 or a compound selected from ~~N-naphthalen-1-yl-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy] acetamide, N-[4-(phenyloxy)phenyl] 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy] acetamide, N-(3,4-dimethylphenyl) 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy] acetamide, N-(2,3-dimethylphenyl) 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy] acetamide, N-(2,4-dimethylphenyl) 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy] acetamide, N-(2,5-dimethylphenyl) 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy] acetamide, N-(3,5-dimethylphenyl) 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy] acetamide, N-(2,6-dimethylphenyl) 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy] acetamide, 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy] N-(2,4,6-trimethylphenyl)acetamide, N-(2-ethylphenyl) 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy] acetamide, N-(4-ethylphenyl) 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy] acetamide, N-(2,6-diethylphenyl) 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy] acetamide, N-[2-(methyloxy)phenyl] 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy] acetamide, N-[2-(ethyloxy)phenyl] 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy] acetamide, N-[3-(ethyloxy)phenyl] 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy] acetamide, N-[2,4-bis(methyloxy)phenyl] 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy] acetamide, N-[4-(dimethylamino)phenyl] 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy] acetamide, N-(2,3-~~

dichlorophenyl) 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy] acetamide, N-(4-chloro-3-methylphenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide, N-(4-bromophenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide, N-(2-fluorophenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide, N-(4-fluorophenyl)-2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide, 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]N-[2-(trifluoromethyl)phenyl]acetamide, 2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]N-[3-(trifluoromethyl)phenyl]acetamide, methyl 4-[[[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetyl]amino] benzoate, ethyl 4-[[[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetyl]amino] benzoate, 3-[[[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetyl]amino] benzoic acid, N-[3-(methyloxy)phenyl]2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide, N-[4-(methyloxy)phenyl]2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide, N-[2-chloro-5-(trifluoromethyl)phenyl]2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide, N-[4-chloro-3-(trifluoromethyl)phenyl]2-[[3-(4H-1,2,4-triazol-4-yl)phenyl]oxy]acetamide, N-(4-chlorophenyl)2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide, N-(4-aminophenyl)2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide, and N-(4-acetylphenyl)2-[[3-(1H-tetrazol-1-yl)phenyl]oxy]acetamide, to a cell or a plurality of cells.